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NEWS	1		Web Page URLs for STN Seminar Schedule - N. America
NEWS	2		"Ask CAS" for self-help around the clock
NEWS	3	JAN 17	Pre-1988 INPI data added to MARPAT
NEWS	4	FEB 21	STN AnaVist, Version 1.1, lets you share your STN AnaVist visualization results
NEWS	5	FEB 22	The IPC thesaurus added to additional patent databases on STN
NEWS	6	FEB 22	Updates in EPFULL; IPC 8 enhancements added
NEWS	7	FEB 27	New STN AnaVist pricing effective March 1, 2006
NEWS	8	MAR 03	Updates in PATDPA; addition of IPC 8 data without attributes
NEWS	9	MAR 22	EMBASE is now updated on a daily basis
NEWS	10	APR 03	New IPC 8 fields and IPC thesaurus added to PATDPAFULL
NEWS	11	APR 03	Bibliographic data updates resume; new IPC 8 fields and IPC thesaurus added in PCTFULL
NEWS	12	APR 04	STN AnaVist \$500 visualization usage credit offered
NEWS	13	APR 12	LINSPEC, learning database for INSPEC, reloaded and enhanced
NEWS	14	APR 12	Improved structure highlighting in FQHIT and QHIT display in MARPAT
NEWS	15	APR 12	Derwent World Patents Index to be reloaded and enhanced during second quarter; strategies may be affected
NEWS	16	MAY 10	CA/CAPLUS enhanced with 1900-1906 U.S. patent records
NEWS	17	MAY 11	KOREAPAT updates resume
NEWS	18	MAY 19	Derwent World Patents Index to be reloaded and enhanced
NEWS	19	MAY 30	IPC 8 Rolled-up Core codes added to CA/CAPLUS and USPATFULL/USPAT2
NEWS	20	MAY 30	The F-Term thesaurus is now available in CA/CAPLUS
NEWS	21	JUN 02	The first reclassification of IPC codes now complete in INPADOC
NEWS EXPRESS		JUNE 16	CURRENT WINDOWS VERSION IS V8.01b, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 23 MAY 2006.
NEWS HOURS			STN Operating Hours Plus Help Desk Availability
NEWS LOGIN			Welcome Banner and News Items
NEWS IPC8			For general information regarding STN implementation of IPC 8
NEWS X25			X.25 communication option no longer available after June 2006

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 13:32:32 ON 19 JUN 2006

=> FIL STNGUIDE

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'STNGUIDE' ENTERED AT 13:32:39 ON 19 JUN 2006

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FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Jun 16, 2006 (20060616/UP).

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.06	0.27

FILE 'REGISTRY' ENTERED AT 13:32:56 ON 19 JUN 2006

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STRUCTURE FILE UPDATES: 18 JUN 2006 HIGHEST RN 888212-64-4

DICTIONARY FILE UPDATES: 18 JUN 2006 HIGHEST RN 888212-64-4

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TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

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\*\*\*\*\*  
 \*  
 \* The CA roles and document type information have been removed from \*  
 \* the IDE default display format and the ED field has been added, \*  
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 \* available and contains the CA role and document type information. \*  
 \*  
 \*\*\*\*\*

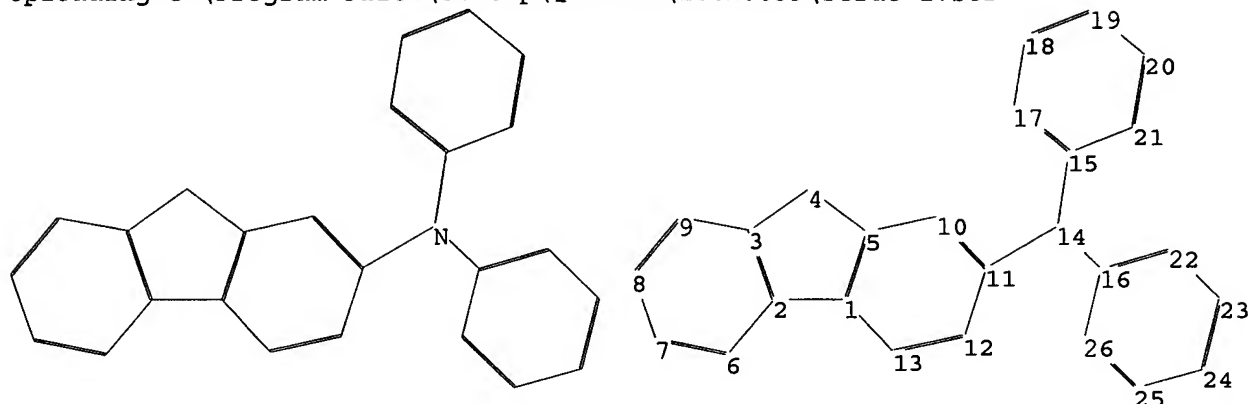
Structure search iteration limits have been increased. See HELP SLIMITS for details.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10620839\Struc 1.str



chain nodes :

14

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 15 16 17 18 19 20 21 22 23 24  
25 26

chain bonds :

11-14 14-15 14-16

ring bonds :

1-2 1-5 1-13 2-3 2-6 3-4 3-9 4-5 5-10 6-7 7-8 8-9 10-11 11-12 12-13  
15-17 15-21 16-22 16-26 17-18 18-19 19-20 20-21 22-23 23-24 24-25 25-26

exact/norm bonds :

1-2 3-4 4-5 11-14 14-15 14-16

normalized bonds :

1-5 1-13 2-3 2-6 3-9 5-10 6-7 7-8 8-9 10-11 11-12 12-13 15-17 15-21  
16-22 16-26 17-18 18-19 19-20 20-21 22-23 23-24 24-25 25-26

Match level :

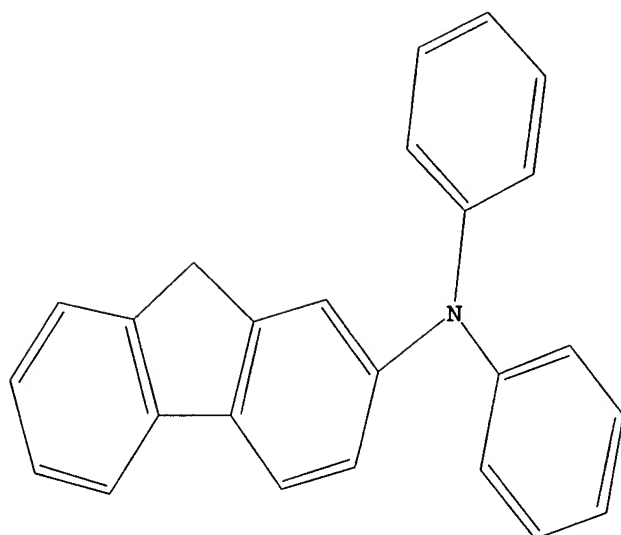
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11:Atom 12:Atom 13:Atom 14:CLASS 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom  
20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom

L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> l1

SAMPLE SEARCH INITIATED 13:33:53 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 3459 TO ITERATE

57.8% PROCESSED 2000 ITERATIONS

50 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

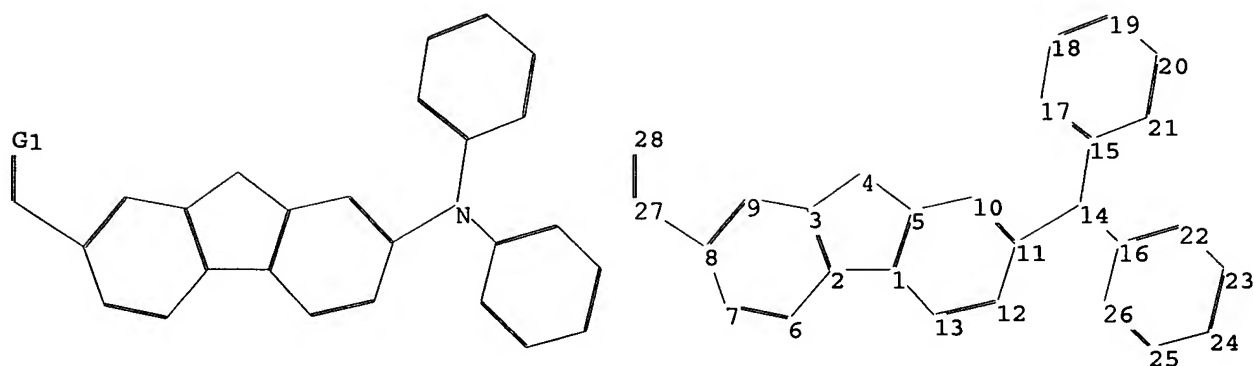
PROJECTED ITERATIONS: 65653 TO 72707

PROJECTED ANSWERS: 1971 TO 3355

L2 50 SEA SSS SAM L1

=>

Uploading C:\Program Files\Stnexp\Queries\10620839\Struc 2.str



```

chain nodes :
14 27 28
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 15 16 17 18 19 20 21 22 23 24
25 26
chain bonds :
8-27 11-14 14-15 14-16 27-28
ring bonds :
1-2 1-5 1-13 2-3 2-6 3-4 3-9 4-5 5-10 6-7 7-8 8-9 10-11 11-12 12-13
15-17 15-21 16-22 16-26 17-18 18-19 19-20 20-21 22-23 23-24 24-25 25-26
exact/norm bonds :
1-2 3-4 4-5 11-14 14-15 14-16 27-28
exact bonds :
8-27
normalized bonds :
1-5 1-13 2-3 2-6 3-9 5-10 6-7 7-8 8-9 10-11 11-12 12-13 15-17 15-21
16-22 16-26 17-18 18-19 19-20 20-21 22-23 23-24 24-25 25-26

```

G1:C,O,N

Match level :

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1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:CLASS 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:CLASS 28:CLASS

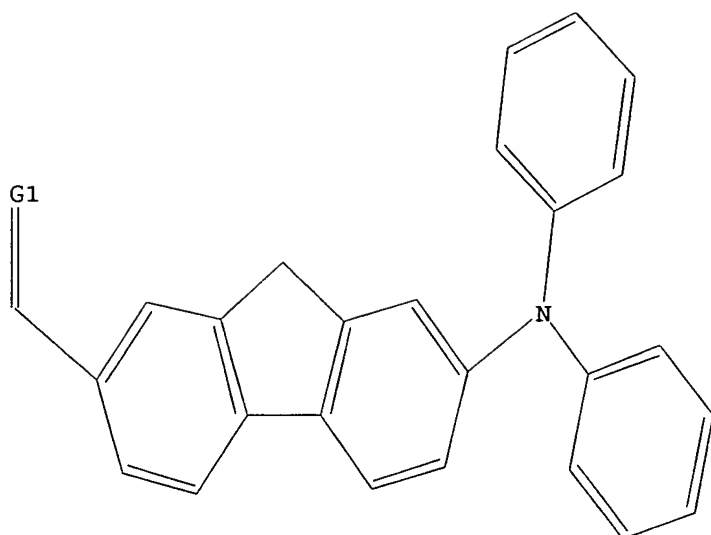
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L3 STRUCTURE UPLOADED

=> d

L3 HAS NO ANSWERS

L3 STR



G1 C,O,N

Structure attributes must be viewed using STN Express query preparation.

=> l3

SAMPLE SEARCH INITIATED 13:35:30 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 2209 TO ITERATE

90.5% PROCESSED 2000 ITERATIONS 1 ANSWERS  
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 41361 TO 46999  
PROJECTED ANSWERS: 1 TO 85

L4 1 SEA SSS SAM L3

=> l3 full

FULL SEARCH INITIATED 13:35:34 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 43292 TO ITERATE

100.0% PROCESSED 43292 ITERATIONS 74 ANSWERS  
SEARCH TIME: 00.00.01

L5 74 SEA SSS FUL L3

=> file medline caplus  
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
168.70	168.97

FULL ESTIMATED COST

FILE 'MEDLINE' ENTERED AT 13:35:43 ON 19 JUN 2006

FILE 'CAPLUS' ENTERED AT 13:35:43 ON 19 JUN 2006

10620839.trn

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=> 15

L6 68 L5

=> d ibib abs hitstr 1-68

L6 ANSWER 1 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2006:290200 CAPLUS

TITLE: Synthesis and characterization of photoresponsive  
diphenylaminofluorene chromophore adducts of  
[60]fullerene

AUTHOR(S): Padmawar, Prashant A.; Canteenwala, Taizoon; Verma,  
Sarika; Tan, Loon-Seng; Chiang, Long Y.

CORPORATE SOURCE: Department of Chemistry, University of Massachusetts  
Lowell, Lowell, 01854, USA

SOURCE: Journal of Materials Chemistry (2006), 16(14),  
1366-1378

CODEN: JMACEP; ISSN: 0959-9428

PUBLISHER: Royal Society of Chemistry

DOCUMENT TYPE: Journal

LANGUAGE: English

AB A class of acceptor-keto-donor structures as hindered 9,9-di(3,5,5-trimethylhexyl)-2-diphenylaminofluoreno-methano[60]fullerene C60(>DPAF-C9) and the related bis-adducts C60(>DPAF-C9)2 and C60(>DPAF-C2)2 were synthesized. They are derivs. of multiphoton absorptive C60(>DPAF-C2) showing enhanced cross-sections of simultaneous two-photon absorption under laser excitation at 800 nm in nanosecond region. Mol. synthesis of these C60-DPAF conjugates involved the covalent attachment of a diphenylaminofluorene moiety to methano[60]fullerene via a keto linkage for increasing mol. acceptor-donor polarization of the chromophore in conjunction with the fullerene cage. Preparation of 7-(1,2-dihydro-1,2-methanofullerene[60]-61-carbonyl)-9,9-dialkyl-2-diphenylaminofluorene C60(>DPAF-Cn) involved cyclopropanation of C60 with a key synthon 7- $\alpha$ -bromoacetyl-9,9-dialkyl-2-diphenylaminofluorene. Synthesis of this synthon was achieved by a three-steps procedure starting from 2-bromofluorene via dialkylation at C9 of the fluorene ring, attachment of a diphenylamino group at C2 of dialkylfluorene, and Friedel-Craft acylation of the  $\alpha$ -bromoacetyl group at C7 of diphenylaminofluorene. All C60-DPAF derivs. were fully characterized with the chemical structures confirmed by various spectroscopic analyses and validated by the single-crystal structural anal. data of C60(>DPAF-C2). Strong solvent-sensitive fluorescence quenching phenomena of C60(>DPAF-C2), C60(>DPAF-C9), and C60(>DPAF-C9)2 were noticed, showing no fluorescence band above 700 nm in more polar solvents, such as DMF, PhCN, and THF, while in less polar solvents (toluene, CHCl<sub>3</sub>, and CS<sub>2</sub>) a fullerenyl fluorescence band at 700-710 nm was observed. It was attributed to the occurrence of electron transfer via the singlet excited state of the fullerene moiety 1C60\*(>DPAF-Cn) in the former group of the solvents. On the contrary, energy transfer processes from DPAF-Cn moiety to the fullerene cage are favored in the latter group of the solvents.

IT 486998-58-7P

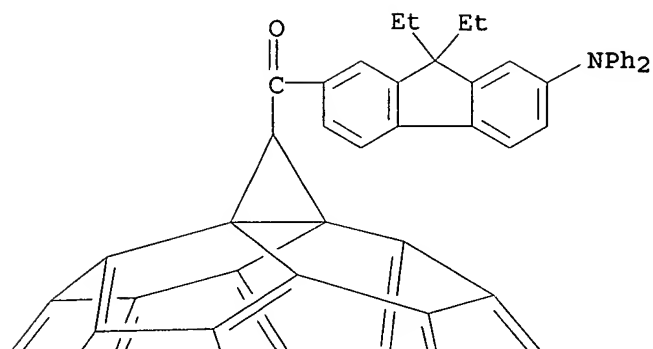
RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

(monoadduct; synthesis and structural characterization and photophys. properties of mono- and bis-adducts of diphenylaminofluorene with methanofullerene)

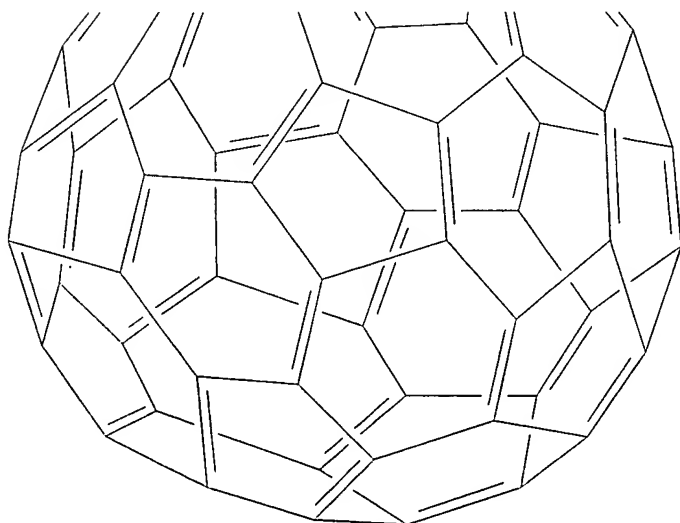
RN 486998-58-7 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

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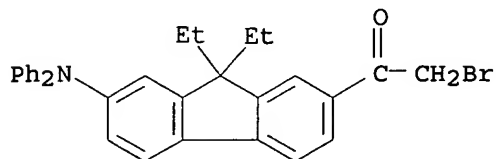
IT 486998-57-6P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)



(reaction with C60 fullerene in presence of DBU)

RN 486998-57-6 CAPLUS

CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI)  
(CA INDEX NAME)

REFERENCE COUNT: 59 THERE ARE 59 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 2 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2006:161203 CAPLUS

DOCUMENT NUMBER: 144:243478

TITLE: Photosensitive polymeric material for worm optical data storage with two-photon fluorescent readout

INVENTOR(S): Belfield, Kevin D.

PATENT ASSIGNEE(S): University of Central Florida Research Foundation, Inc., USA

SOURCE: U.S., 33 pp.  
CODEN: USXXAM

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 7001708	B1	20060221	US 2002-306960	20021127
PRIORITY APPLN. INFO.:			US 2001-333972P	P 20011128
			US 2001-339283P	P 20011211

AB Image formation via photoinduced fluorescence changes in a polymeric medium with two-photon fluorescence readout of a multi-layer structure. Fluorophore-containing polymers, possessing one or more basic functional groups, underwent protonation in the presence of a photoinduced acid generator upon exposure to a broad-band UV light source or fast-pulsed red to near-IR laser irradiation. Solution studies demonstrated formation of monoprotonated and diprotonated species upon irradiation, each resulting in distinctly different absorption and fluorescence properties. The fluorescence of the original, neutral, fluorophore was reduced upon monoprotonation, leading to a concomitant increase in fluorescence at longer wavelengths due to the monoprotonated form, the basis for multichannel data readout. Expts. in polymer films demonstrate the changes in fluorescence properties of the photosensitive polymer compns. and polymers can be employed for a high storage d., write-once read-many (WORM) data storage medium with two-photon fluorescence readout. Two-channel, two-photon fluorescence imaging provided both "pos." and "neg." image readout capability.

IT 745079-44-1P

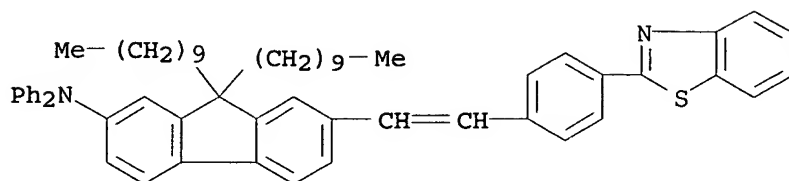
RL: SPN (Synthetic preparation); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)

(photosensitive polymeric material for WORM optical data storage)

RN 745079-44-1 CAPLUS

CN 9H-Fluoren-2-amine, 7-[2-[4-(2-benzothiazolyl)phenyl]ethenyl]-9,9-didecyl-

N,N-diphenyl- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 38 THERE ARE 38 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 3 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:1334086 CAPLUS

DOCUMENT NUMBER: 144:149941

TITLE: Photoinduced Processes in a Tricomponent Molecule Consisting of Diphenylaminofluorene-Dicyanoethylene-Methano[60]fullerene

AUTHOR(S): El-Khouly, Mohamed E.; Padmawar, Prashant; Araki, Yasuyuki; Verma, Sarika; Chiang, Long Y.; Ito, Osamu  
CORPORATE SOURCE: Institute of Multidisciplinary Research for Advanced Materials, Tohoku University, Sendai, 980-8577, Japan  
SOURCE: Journal of Physical Chemistry A (2006), 110(3), 884-891

CODEN: JPCAFH; ISSN: 1089-5639

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

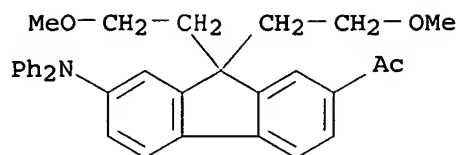
AB Photoinduced intramol. processes in a tricomponent mol. C60(>(CN)2-DPAF), consisting of an electron-accepting methano[60]fullerene moiety (C60>) covalently bound to an electron-donating diphenylaminofluorene (DPAF) unit via a bridging dicyanoethylenyl group [(CN)2], were investigated in comparison with (CN)2-DPAF. On the basis of the MO calcns., the lowest charge-separated state of C60(>(CN)2-DPAF) is suggested to be C60•-(>(CN)2-DPAF•+) with the neg. charge localized on the fullerene cage, while the upper state is C60(>(CN)2•-DPAF•+). The excited-state events of C60(>(CN)2-DPAF) were monitored by both time-resolved emission and nanosecond transient absorption techniques. In both nonpolar and polar solvents, the excited charge-transfer state decayed mainly through initial energy-transfer process to the C60 moiety yielding the corresponding 1C60\*, from which charge separation took place leading to the formation of C60•-(>(CN)2-DPAF•+) in a fast rate and high efficiency. In addition, multistep charge separation from C60(>(CN)2•-DPAF•+) to C60•-(>(CN)2-DPAF•+) may be possible with the excitation of charge-transfer band. The lifetimes of C60•-(>(CN)2-DPAF•+) are longer than the previously reported methano[60]fullerene-diphenylaminofluorene C60(>(C:O)-DPAF) with the C60 and DPAF moieties linked by a methanoketo group. These findings suggest an important role of dicyanoethylenyl group as an electron mediating bridge in C60(>(CN)2-DPAF).

IT 874200-94-9

RL: PRP (Properties); RCT (Reactant); RACT (Reactant or reagent)  
(condensation with malononitrile; photoinduced processes in a (diphenylamino)fluorene-dicyanoethylene-methano[60]fullerene triad)

RN 874200-94-9 CAPLUS

CN Ethanone, 1-[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



IT 649724-49-2

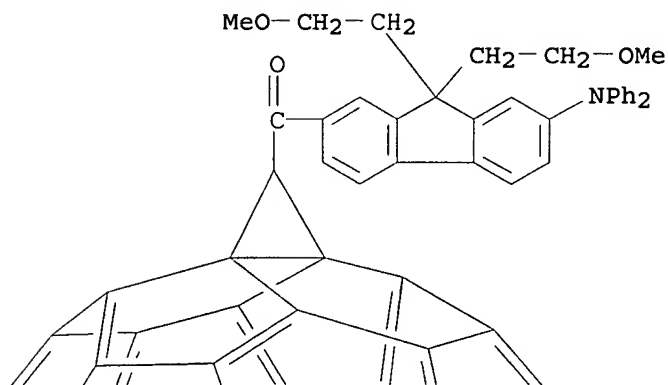
RL: RCT (Reactant); RACT (Reactant or reagent)

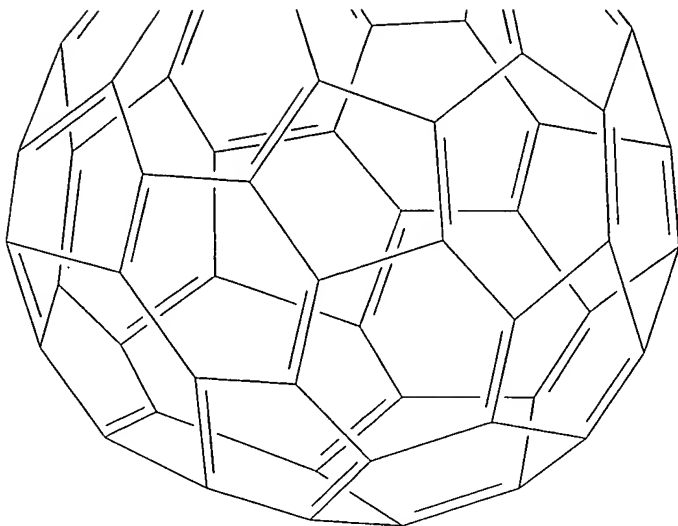
(condensation with malononitrile; photoinduced processes in a  
(diphenylamino)fluorene-dicyanoethylene-methano[60]fullerene triad)

RN 649724-49-2 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-Ih-3'-yl[7-(diphenylamino)-  
9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

PAGE 1-A





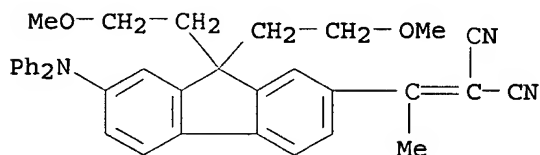
IT 874200-97-2

RL: FMU (Formation, unclassified); PRP (Properties); FORM (Formation, nonpreparative)

(formation by chemical oxidation; photoinduced processes in a (diphenylamino)fluorene-dicyanoethylene-methano[60]fullerene triad)

RN 874200-97-2 CAPLUS

CN Propanedinitrile, [1-[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]ethylidene]-, radical ion(1+) (9CI) (CA INDEX NAME)



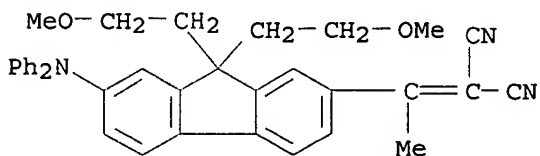
IT 874200-95-0P

RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

(reference compound; photoinduced processes in a (diphenylamino)fluorene-dicyanoethylene-methano[60]fullerene triad)

RN 874200-95-0 CAPLUS

CN Propanedinitrile, [1-[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]ethylidene]- (9CI) (CA INDEX NAME)



IT 874200-96-1P

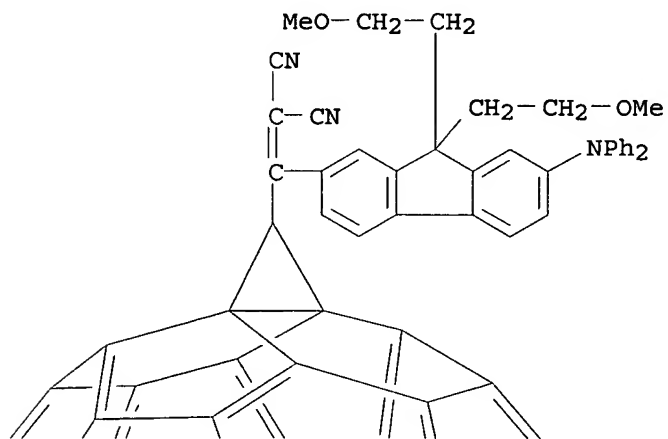
RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

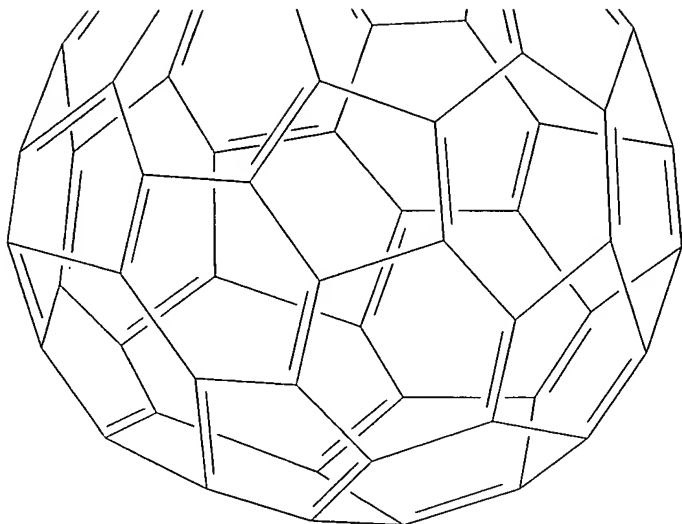
(target triad; photoinduced processes in a (diphenylamino)fluorene-dicyanoethylene-methano[60]fullerene triad)

RN 874200-96-1 CAPLUS

CN Propanedinitrile, [3'H-cyclopropa[1,9][5,6]fulleren-C60-Ih-3'-yl[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]methylene] - (9CI)  
(CA INDEX NAME)

PAGE 1-A





REFERENCE COUNT: 90 THERE ARE 90 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 4 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:1323366 CAPLUS

DOCUMENT NUMBER: 144:483596

TITLE: Singlet oxygen production by amphiphilic C60 derivatives and its correlation to cell cytotoxicity in vitro

AUTHOR(S): So, Grace; Karotki, Aliaksandr; Verma, Sarika; Hauck, Tanya S.; Wilson, Brian; Pritzker, Kenneth P. H.; Chiang, Long

CORPORATE SOURCE: Department of Materials Science and Engineering, Univ. of Toronto, Can.

SOURCE: Proceedings of SPIE-The International Society for Optical Engineering (2005), 5969(Photonic Applications in Biosensing and Imaging), 59690D/1-59690D/8  
CODEN: PSISDG; ISSN: 0277-786X

PUBLISHER: SPIE-The International Society for Optical Engineering

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Fullerene derivs. have appealing properties that can potentially be used in materials science and medical applications. In particular, fullerenes are known to produce reactive oxygen species upon their excitation with light. This makes them particularly attractive as photosensitizers for photodynamic therapy (PDT). Photodynamic therapy is a new modality of treatment of cancer as well as some non-cancerous conditions. It involves the combined actions of a drug (photosensitizer) and light to produce a cytotoxic effect. Water-soluble hexa(sulfo-n-butyl)[60]fullerenes (FC4S) was reported recently to generate singlet oxygen ( $^1O_2$ ) and superoxide radical ( $O_2^{\cdot-}$ ) upon its excitation with light, making it a promising candidate for PDT treatments. Recently, we synthesized new amphiphilic fullerene derivs., namely, [60]fullerene-diphenylaminofluorene-oligo(ethylene glycol) conjugates, C60(>DPAF-PEG600) and C60(>DPAF-PEG2000), as potential photosensitizers. In this paper we compare FC4S to PEG-based fullerenes in terms of their singlet oxygen photosensitization ability. We measured time-resolved kinetics of singlet

oxygen luminescence photosensitized by excitation of fullerenes via a 10 ns pulsed laser at 523 nm. For FC4S we observed "normal" kinetics with a monoexponential decay profile giving a time constant 3.8 us in water. In contrast, for the case of C60(>DPAF-PEG600) and C60(>DPAF-PEG2000), a non-monoexponential decay profile with a long tail (.apprx. 102 μs) in water was observed. We hypothesize that this is due to formation of vesicles by PEG fullerenes in aqueous solution. To investigate photodynamic activity of these fullerene derivs. in vitro, we used HeLa human adenocarcinoma and B16 mouse melanoma cell lines. FC4S showed clear photodynamic effects in both cell lines. The total fluence required to kill 50% of the cells at the drug concentration of 20 μM was 36 Jcm<sup>-2</sup> for HeLa cells and 72 Jcm<sup>-2</sup> for B16 cells. Neither PEG-based fullerene derivs. showed any appreciable photodynamic activity, possibly, due to low efficiency of singlet oxygen generation.

IT 851204-04-1

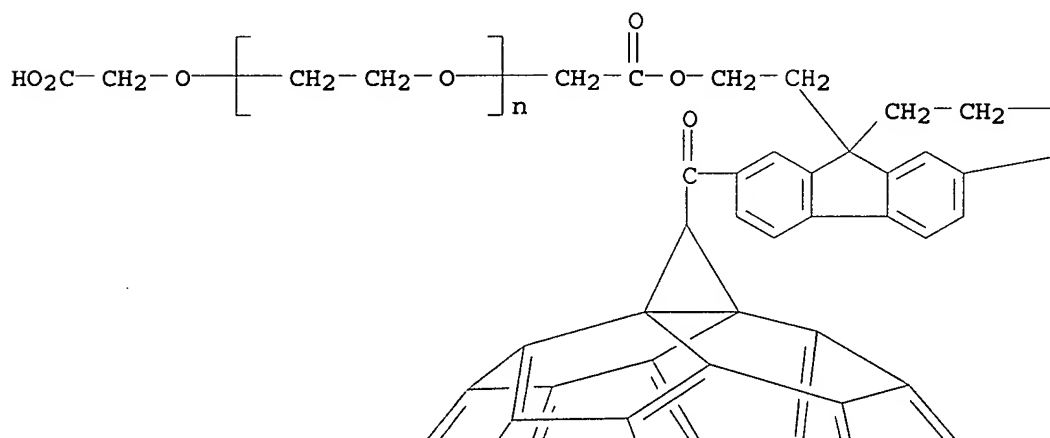
RL: DMA (Drug mechanism of action); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

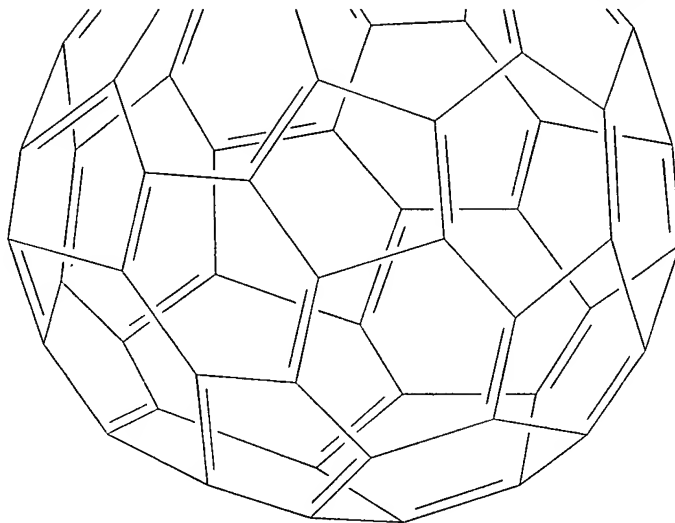
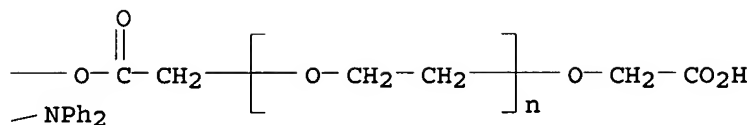
(singlet oxygen production by amphiphilic C60 derivs. and its correlation to cell cytotoxicity in vitro)

RN 851204-04-1 CAPLUS

CN Poly(oxy-1,2-ethanediyl), α,α'-[[2-(3'H-cyclopropa[1,9][5,6]fullerene-C60-1h-3'-ylcarbonyl)-7-(dipropylamino)-9H-fluoren-9-ylidene]bis[2,1-ethanediyl]oxy(2-oxo-2,1-ethanediyl)]]bis[ω-(carboxymethoxy)-(9CI) (CA INDEX NAME)

PAGE 1-A





REFERENCE COUNT: 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 5 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2005:1228344 CAPLUS  
 DOCUMENT NUMBER: 144:129366  
 TITLE: Synthesis of Water-Soluble Highly Two-Photon Responsive [60]Fullerene-Diphenylaminofluorene Chromophore Dyads  
 AUTHOR(S): Verma, Sarika; Padmawar, Prashant; Hauck, Tanya; Canteenwala, Taizoon; Chiang, Long; Pritzker, Kenneth  
 CORPORATE SOURCE: Department of Chemistry, University of Massachusetts, Lowell, MA, USA  
 SOURCE: Journal of Macromolecular Science, Part A: Pure and Applied Chemistry (2005), A42(11), 1497-1505  
 CODEN: JSPCE6; ISSN: 1060-1325  
 PUBLISHER: Taylor & Francis, Inc.  
 DOCUMENT TYPE: Journal



LANGUAGE: English

AB We have synthesized an amphiphilic donor-acceptor type [60]fullerene-diphenylaminofluorene (DPAF) containing oligo(ethylene glycol) moieties, denoted C60-DPAF-EGn. The hydrophobic C60-DPAF moiety consisting of an electron-donating diphenylaminofluorene unit and an electron-accepting C60 cage shows large cross-sections of two-photon absorptivity (2PA) in the nanosecond region. The attachment of two hydrophilic oligo(ethylene glycol) segments on DPAF mainly increases water-solubility of the corresponding chromophoric macromols. that enhances their potential application in biomedical treatments. Here, we present the synthesis and self-assembly study of these amphiphilic materials in aqueous solution New C60-DPAF-EGn compds. were structurally characterized by various spectroscopic methods including, <sup>1</sup>H-NMR, <sup>13</sup>C-NMR, FT-IR, and UV spectroscopies. In the vesicle formation investigation, UV-visible absorption spectra, dynamic light scattering measurements, and transmission electron microscopy were used as primary methods for morphol. characterization of mol. self-assembly behavior of these amphiphilic [60]fullerene-diphenylaminofluorene chromophore dyads.

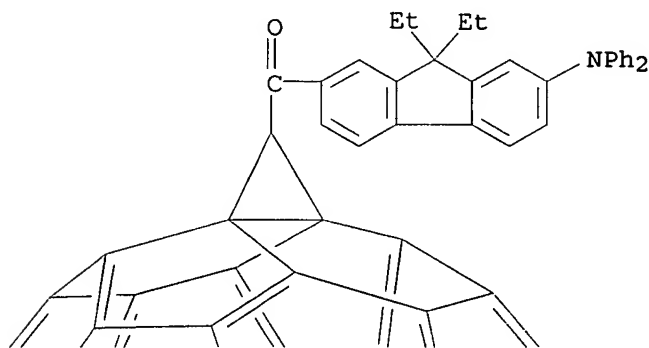
IT 486998-58-7 649724-49-2 649724-50-5

RL: RCT (Reactant); RACT (Reactant or reagent)  
(in synthesis of water-soluble highly two-photon responsive  
[60]Fullerene-diphenylaminofluorene chromophore dyads)

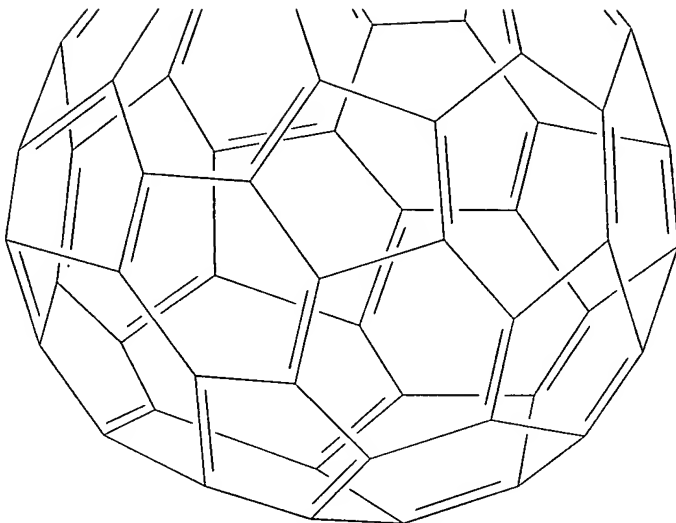
RN 486998-58-7 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

PAGE 1-A

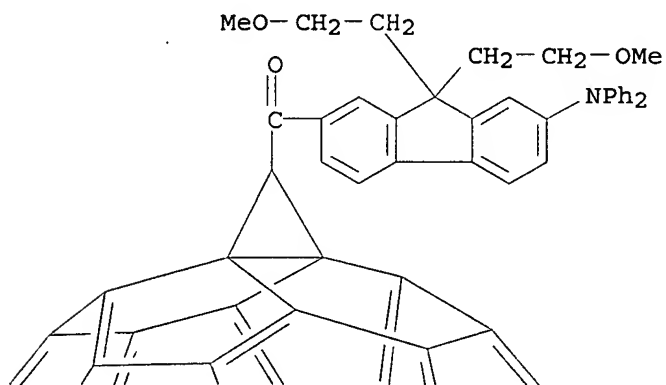


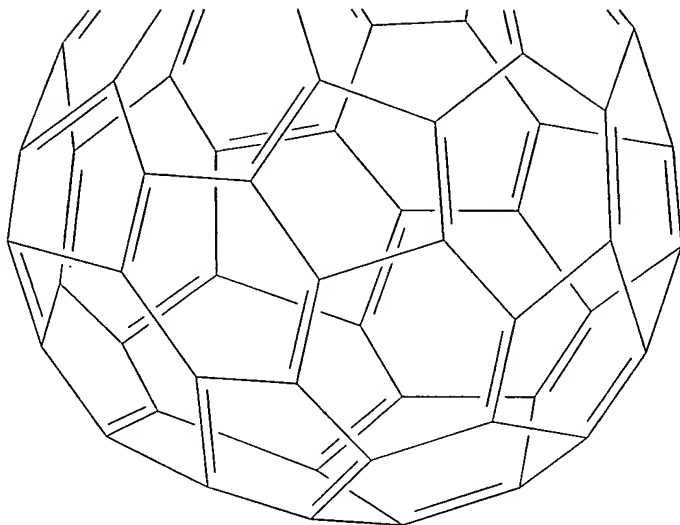
PAGE 2-A



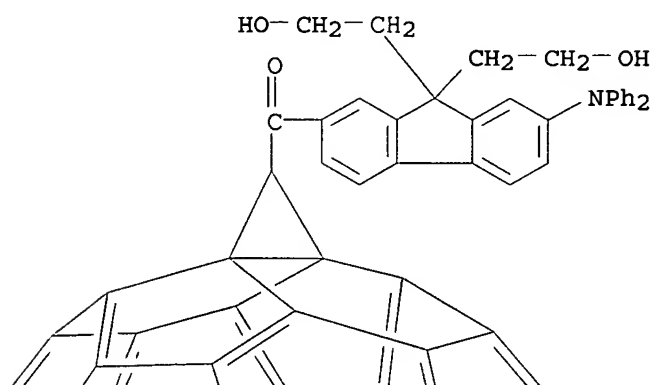
RN 649724-49-2 CAPLUS  
 CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]-(9CI) (CA INDEX NAME)

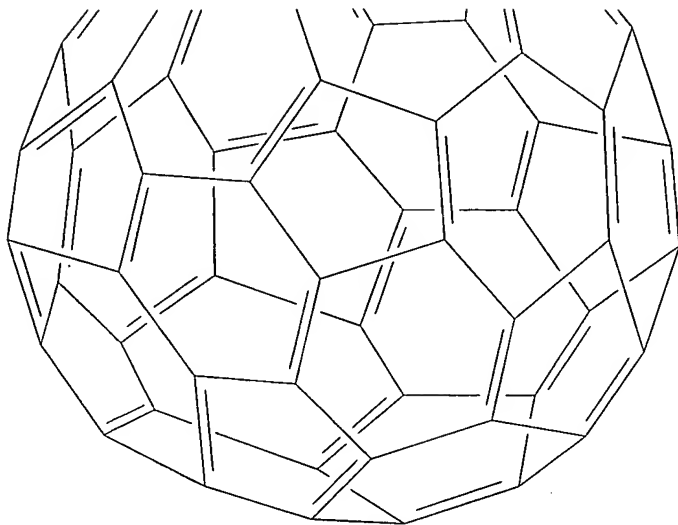
PAGE 1-A





RN 649724-50-5 CAPLUS  
 CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-bis(2-hydroxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



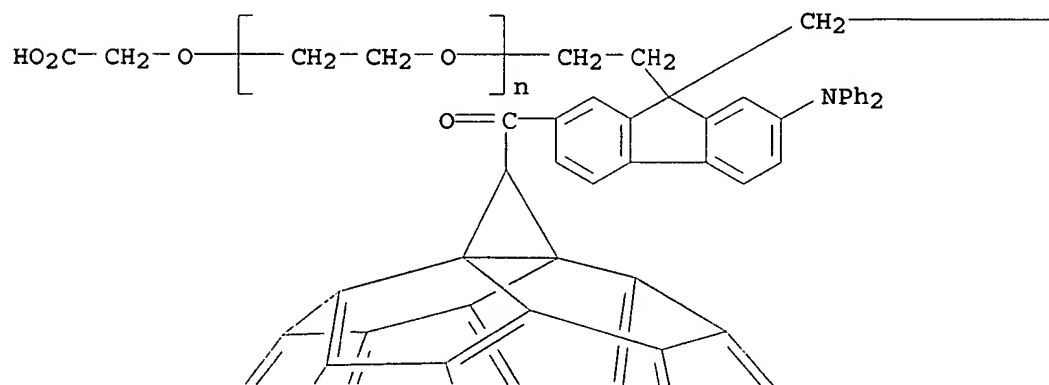


IT 873437-03-7P

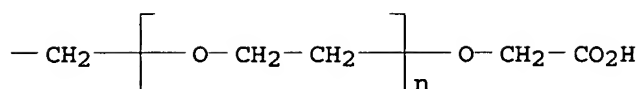
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
 (synthesis of water-soluble highly two-photon responsive  
 [60]Fullerene-diphenylaminofluorene chromophore dyads)

RN 873437-03-7 CAPLUS

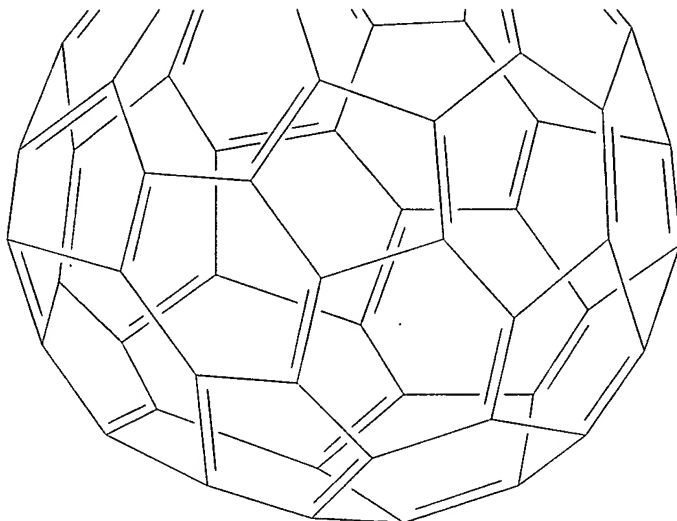
CN Poly(oxy-1,2-ethanediyl),  $\alpha, \alpha'$ -[[2-(3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-ylcarbonyl)-7-(diphenylamino)-9H-fluoren-9-ylidene]di-2,1-ethanediyl]bis[ $\omega$ -(carboxymethoxy)- (9CI)  
 (CA INDEX NAME)



PAGE 1-B



PAGE 2-A



REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 6 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2005:1085998 CAPLUS  
 DOCUMENT NUMBER: 144:42395  
 TITLE:  $\pi$ -Conjugated Dendritic Nanosized Chromophore with Enhanced Two-Photon Absorption  
 AUTHOR(S): Zheng, Qingdong; He, Guang S.; Prasad, Paras N.  
 CORPORATE SOURCE: Department of Chemistry, Institute for Lasers, Photonics and Biophotonics, State University of New York at Buffalo, Buffalo, NY, 14260, USA  
 SOURCE: Chemistry of Materials (2005), 17(24), 6004-6011  
 CODEN: CMATEX; ISSN: 0897-4756  
 PUBLISHER: American Chemical Society  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 AB A novel  $\pi$ -conjugated dendritic nanosized chromophore,

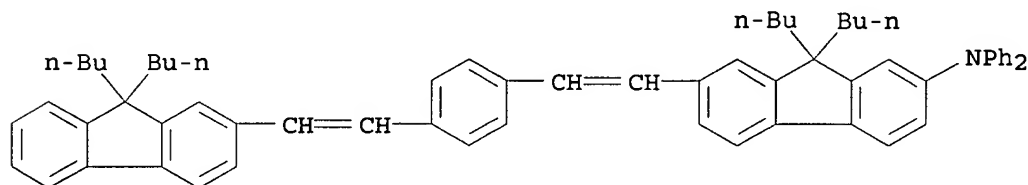
1,3,5-tris(7,12-bis(2-{4-[2-(7-diphenylamino-9,9-dibutyl-9H-fluoren-2-yl)-vinyl]-phenyl}-vinyl)-5,5,10,10,15,15-hexabutyl-10,15-dihydro-5H-diindeno[1,2-a;1',2'-c]fluoren-2-yl)-benzene (3), was synthesized and characterized, together with its corresponding one- and three-branched units. Two-photon absorption (TPA) spectra for these chromophores were measured by the nonlinear transmission spectral technique which utilizes a femtosecond white-light continuum. The TPA peak value for the dendritic chromophore is  $5.47 \times 10^{-20}$  cm<sup>4</sup>/GW, 15.6 times larger than that for the one-branched dipolar chromophore ( $0.35 \times 10^{-20}$  cm<sup>4</sup>/GW), [9,9-dibutyl-7-(2-{4-[2-(9,9-dibutyl-9H-fluoren-2-yl)-vinyl]-phenyl}-vinyl)-9H-fluoren-2-yl]-diphenylamine (1). The three-branched chromophore, 2,7,12-tris(2-{4-[2-(9,9-dibutyl-7-diphenylamino-9H-fluoren-2-yl)-vinyl]-phenyl}-vinyl)-5,5,10,10,15,15-hexabutyl-10,15-dihydro-5H-diindeno[1,2-a;1',2'-c]fluorene (2) has a TPA peak value of  $2.28 \times 10^{-20}$  cm<sup>4</sup>/GW, 6.5 times as large as that for compound 1. The enhanced two-photon absorption found in these two multibranched chromophores is attributed to the fact that the dendritic or the three-branched chromophore has both an extended  $\pi$ -conjugated system and an increased intramol. charge redistribution compared to the one-branched chromophore. With use of comparable structure unit based concns., the two-photon excited (TPE) fluorescence intensity for this dendritic chromophore (0.0033 M solution) was found to be enhanced by a factor of 2.9 compared to that for chromophore 1 (0.02 M solution), which would offer a major advantage in TPE fluorescence related applications.

IT 870772-54-6P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
(one-branched chromophore; synthesis and linear- and two-photon  
absorption and two-photon excited fluorescence of  $\pi$ -conjugated  
dendritic nanosized chromophores containing diphenylamine- and Bu groups)

RN 870772-54-6 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-dibutyl-7-[2-[4-[2-(9,9-dibutyl-9H-fluoren-2-yl)ethenyl]phenyl]ethenyl]-N,N-diphenyl- (9CI) (CA INDEX NAME)

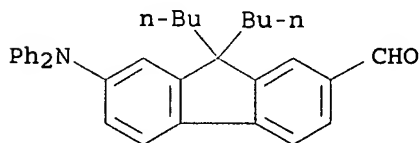


IT 870772-56-8

RL: RCT (Reactant); RACT (Reactant or reagent)  
(synthesis of  $\pi$ -conjugated dendritic nanosized chromophores containing  
diphenylamine- and Bu groups)

RN 870772-56-8 CAPLUS

CN 9H-Fluorene-2-carboxaldehyde, 9,9-dibutyl-7-(diphenylamino)- (9CI) (CA INDEX NAME)



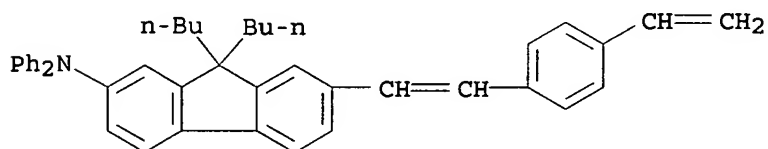
IT 870772-57-9P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(synthesis of  $\pi$ -conjugated dendritic nanosized chromophores containing diphenylamine- and Bu groups)

RN 870772-57-9 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-dibutyl-7-[2-(4-ethenylphenyl)ethenyl]-N,N-diphenyl- (9CI) (CA INDEX NAME)



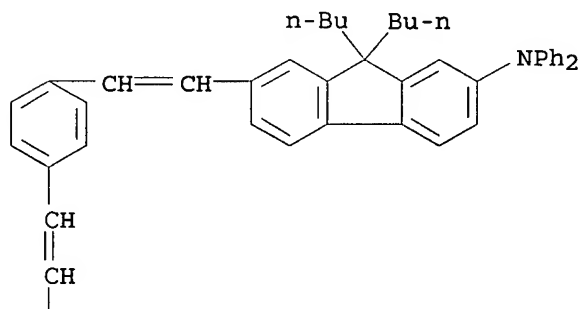
IT 870772-55-7P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)

(three-branched chromophore; synthesis and linear- and two-photon absorption and two-photon excited fluorescence of  $\pi$ -conjugated dendritic nanosized chromophores containing diphenylamine- and Bu groups)

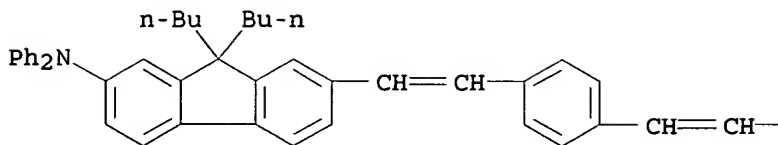
RN 870772-55-7 CAPLUS

CN 9H-Fluoren-2-amine, 7,7',7''-[(5,5,10,10,15,15-hexabutyl-10,15-dihydro-5H-tribenzo[a,f,k]trindene-2,7,12-triyl)tris(2,1-ethenediyl-4,1-phenylene-2,1-ethenediyl)]tris[9,9-dibutyl-N,N-diphenyl- (9CI) (CA INDEX NAME)

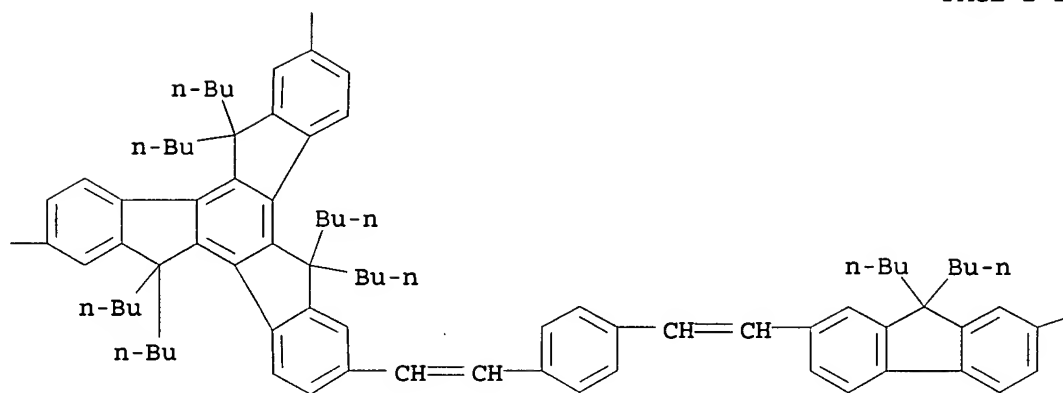


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PAGE 2-B



PAGE 2-C

—NPh<sub>2</sub>

REFERENCE COUNT: 39 THERE ARE 39 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 7 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:1062319 CAPLUS

DOCUMENT NUMBER: 144:23966

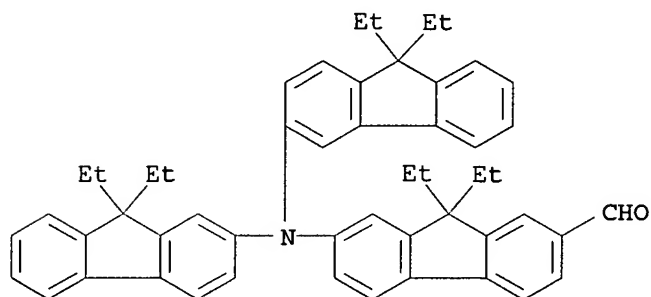
TITLE: New fluorophores based on trifluorenylamine with very large intrinsic three-photon absorption cross sections

AUTHOR(S): Suo, Zhiyong; Drobizhev, Mikhail; Spangler, Charles W.; Christensson, Niklas; Rebane, Alexander

CORPORATE SOURCE: Physics Department and Chemistry and Biochemistry Department, Montana State University, Bozeman, MT,

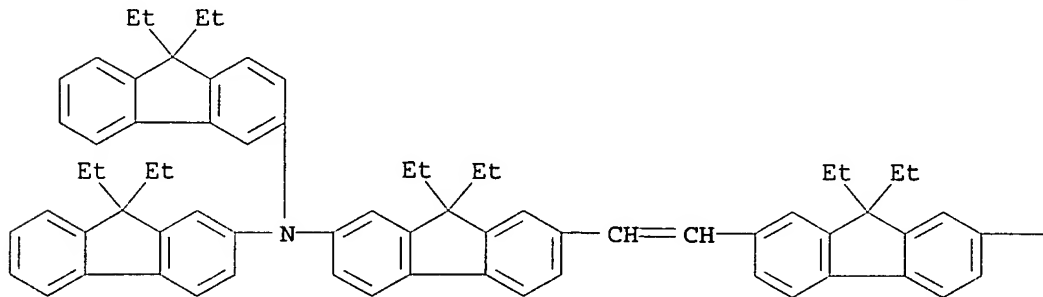


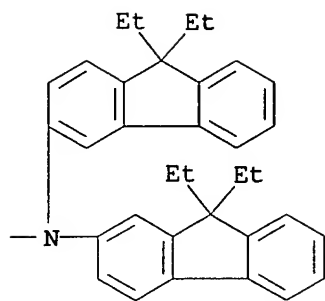
59717, USA  
 SOURCE: Organic Letters (2005), 7(22), 4807-4810  
 CODEN: ORLEF7; ISSN: 1523-7060  
 PUBLISHER: American Chemical Society  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 OTHER SOURCE(S): CASREACT 144:23966  
 AB A new fluorophore, tris(9,9-diethyl-9H-fluorenyl)amine, was synthesized by the Buchwald-Hartwig reaction of 2-aminofluorene, and based on this mol. three more fluorophores were prepared that exhibit a very large intrinsic three-photon absorption in the near-IR region, which scales as a third power of the bridge length.  
 IT 870283-38-8P  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (intermediate; preparation of fluorophores based on trifluorenylamine with very large intrinsic three-photon absorption cross sections)  
 RN 870283-38-8 CAPLUS  
 CN 9H-Fluorene-2-carboxaldehyde, 7-[(9,9-diethyl-9H-fluoren-2-yl)(9,9-diethyl-9H-fluoren-3-yl)amino]-9,9-diethyl- (9CI) (CA INDEX NAME)



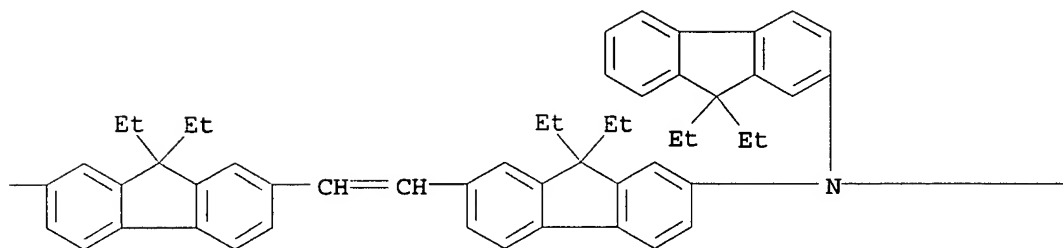
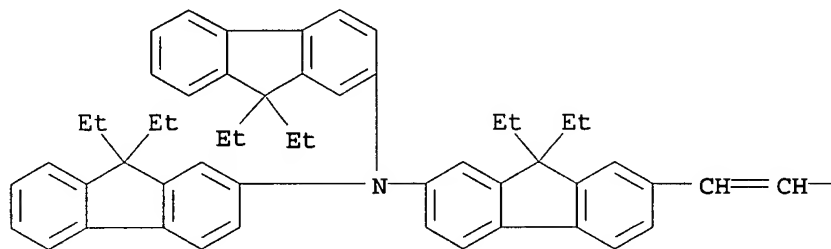
IT 870283-42-4P 870283-43-5P  
 RL: PRP (Properties); SPN (Synthetic preparation); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)  
 (orange fluorophore; preparation of fluorophores based on trifluorenylamine with very large intrinsic three-photon absorption cross sections)  
 RN 870283-42-4 CAPLUS  
 CN 9H-Fluorene-2-amine, 7,7'-(1,2-ethenediyl)bis[N-(9,9-diethyl-9H-fluoren-2-yl)-N-(9,9-diethyl-9H-fluoren-3-yl)-9,9-diethyl- (9CI) (CA INDEX NAME)

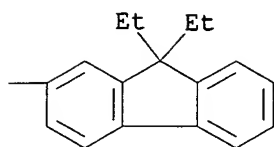
PAGE 1-A





RN 870283-43-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 7,7'-[(9,9-diethyl-9H-fluorene-2,7-diyl)di-2,1-ethenediyl]bis[9,9-diethyl-N,N-bis(9,9-diethyl-9H-fluoren-2-yl) - (9CI)  
 (CA INDEX NAME)





REFERENCE COUNT: 36 THERE ARE 36 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 8 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:988390 CAPLUS

DOCUMENT NUMBER: 143:451502

TITLE: Synthesis of C60-diphenylaminofluorene dyads with two-photon absorbing characteristics

AUTHOR(S): Padmawar, Prashant A.; Canteenwala, Taizoon; Verma, Sarika; Tan, Loon-Seng; He, Guang S.; Prasad, Paras N.; Chiang, Long Y.

CORPORATE SOURCE: Department of Chemistry, Institute of Nanoscience and Engineering Technology, University of Massachusetts, Lowell, MA, 01854, USA

SOURCE: Synthetic Metals (2005), 154(1-3), 185-188

CODEN: SYMEDZ; ISSN: 0379-6779

PUBLISHER: Elsevier B.V.

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Linear A-sp<sup>3</sup>-D mol. conjugates as 9,9-dialkyl-2-diphenylaminofluorene (DPAF-C<sub>2</sub>)-C60 monoadducts were demonstrated using methanoketo unit as a linker to bridge the DPAF donor moiety and the fullerene acceptor chromophore together within a short distance of roughly 2.0 Å. Target products of C60-diphenylaminofluorene dyads C60(>DPAF-C<sub>2</sub>) and C60(>DPAF-C<sub>18</sub>) were synthesized using Bingel cyclopropanation reaction from the corresponding 7-α-bromoacetyl-9,9-dialkyl-2-diphenylaminofluorene precursors. Both dyads C60(>DPAF-C<sub>2</sub>) and C60(>DPAF-C<sub>18</sub>) were characterized by spectroscopic methods and simultaneous two-photon excitation measurements, showing large two-photon absorption cross-sections in the nanosecond regime.

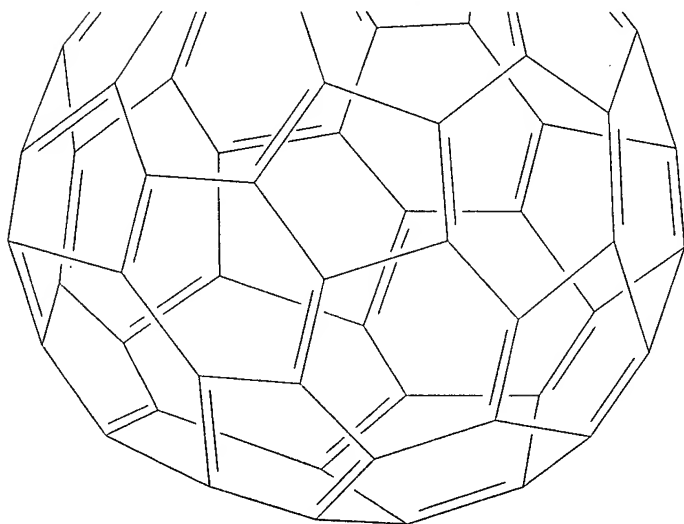
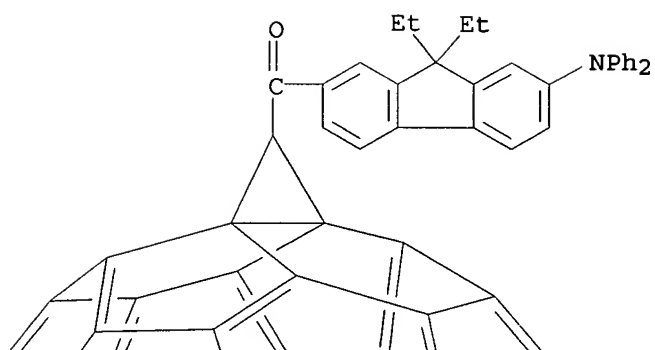
IT 486998-58-7P 868852-15-7P

RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

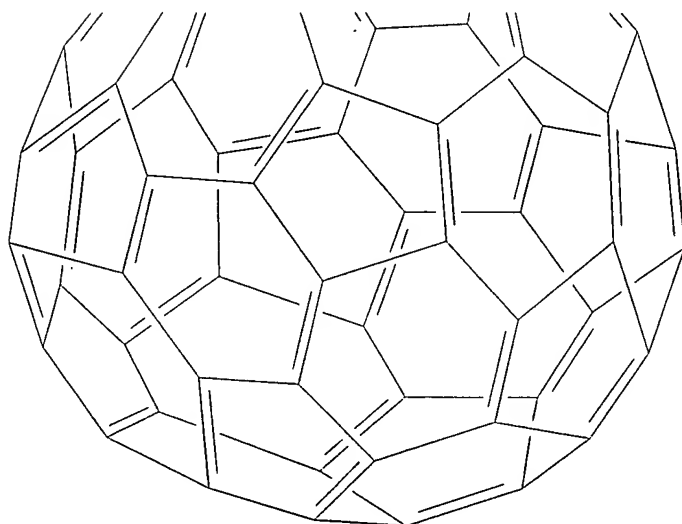
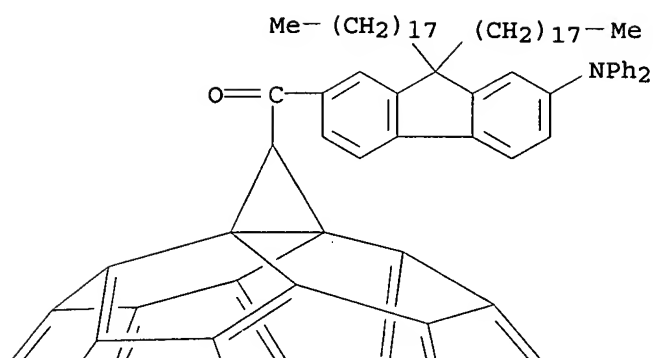
(synthesis and photophys. of C60-diphenylaminofluorene dyads with two-photon absorbing characteristics)

RN 486998-58-7 CAPLUS

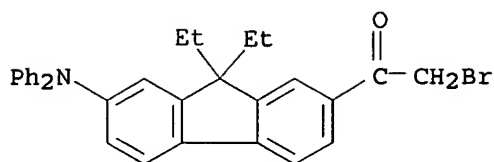
CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



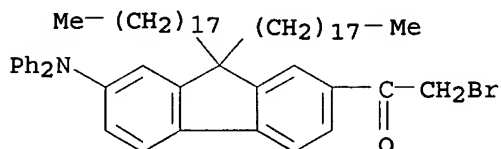
RN 868852-15-7 CAPLUS  
 CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-Ih-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



IT 486998-57-6 868852-16-8  
 RL: RCT (Reactant); RACT (Reactant or reagent)  
 (synthesis of C60-diphenylaminofluorene dyad)  
 RN 486998-57-6 CAPLUS  
 CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI)  
 (CA INDEX NAME)



RN 868852-16-8 CAPLUS  
 CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-dioctadecyl-9H-fluoren-2-yl]-(9CI) (CA INDEX NAME)



REFERENCE COUNT: 20 THERE ARE 20 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 9 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:865315 CAPLUS

DOCUMENT NUMBER: 143:412994

TITLE: Synthesis, two- and three-photon absorption, and optical limiting properties of fluorene-containing ferrocene derivatives

AUTHOR(S): Zheng, Qingdong; He, Guang S.; Lu, Changgui; Prasad, Paras N.

CORPORATE SOURCE: Department of Chemistry, Institute for Lasers, Photonics and Biophotonics, State University of New York at Buffalo, Buffalo, NY, 14260, USA

SOURCE: Journal of Materials Chemistry (2005), 15(34), 3488-3493

CODEN: JMACEP; ISSN: 0959-9428

PUBLISHER: Royal Society of Chemistry

DOCUMENT TYPE: Journal

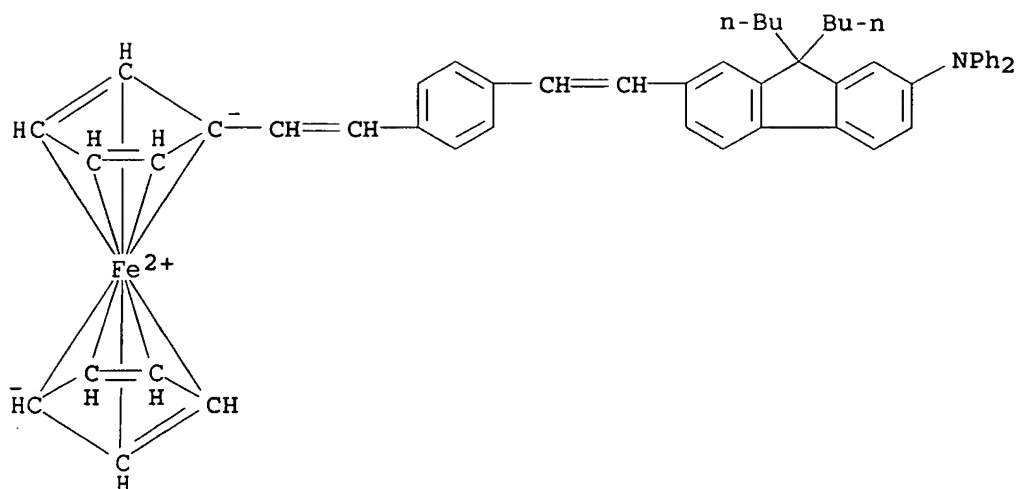
LANGUAGE: English

AB The synthesis and characterization of 2 novel fluorene-containing ferrocene derivs. were reported. The 2-photon absorption spectra and 3-photon absorption cross-section values in the IR region for these 2 chromophores were studied. Together with their thermal stabilities, their linear absorption and emission properties were also studied. The ferrocene derivs. have large 2-photon and 3-photon absorption in IR region as well as excellent thermal stabilities. The 3-photon absorption based optical limiting properties of these 2 ferrocene derivs. were studied by using sub-picosecond IR laser pulses.

IT 867045-90-7P, 9,9-Dibutyl-7-[(E)-2-[4-[(E)-(2-ferrocenylvinyl)]phenyl]vinyl-9H-fluoren-2-yl]diphenylamine  
 RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
 (synthesis, two- and three-photon absorption, and optical limiting properties of fluorene-containing ferrocene derivs.)

RN 867045-90-7 CAPLUS

CN Ferrocene, [(1E)-2-[4-[(1E)-2-[9,9-dibutyl-7-(diphenylamino)-9H-fluoren-2-yl]ethenyl]phenyl]ethenyl]- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 35 THERE ARE 35 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 10 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:780629 CAPLUS

DOCUMENT NUMBER: 144:201429

TITLE: Measurement of the two-photon absorption cross sections of dicyanomethylene-pyrans by the z-scan method

AUTHOR(S): Chunosova, S. S.; Svetlichnyi, V. A.; Meshalkin, Yu. P.

CORPORATE SOURCE: Institute of Laser Physics, Siberian Branch, Russian Academy of Sciences, Novosibirsk, 630090, Russia

SOURCE: Quantum Electronics (2005), 35(5), 415-418  
CODEN: QUELEZ; ISSN: 1063-7818

PUBLISHER: Turpion Ltd.

DOCUMENT TYPE: Journal

LANGUAGE: English

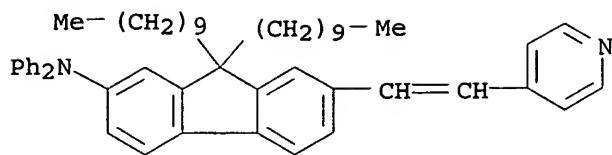
AB The two-photon absorption cross sections of three dicyanomethylene-pyrane DCM dyes excited by a femtosecond pulsed from a Ti:sapphire laser are measured by the open-aperture z-scan method. The exptl. transmission curves are approximated by the corresponding model curves to give the two-photon absorption cross sections equal to  $5.1 \times 10^3$  GM,  $5.7 \times 10^3$  GM, and  $8.4 \times 10^3$  GM for DCM, DCM-17, and DCM-doa, resp. (1 GM =  $10^{-50}$  cm<sup>4</sup> s phot<sup>-1</sup> mol<sup>-1</sup>).

IT 191667-13-7, AF-50

RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); PROC (Process)  
(measurement of two-photon absorption cross sections of dicyanomethylene-pyrans by the z-scan method)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



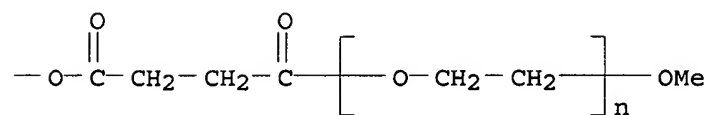
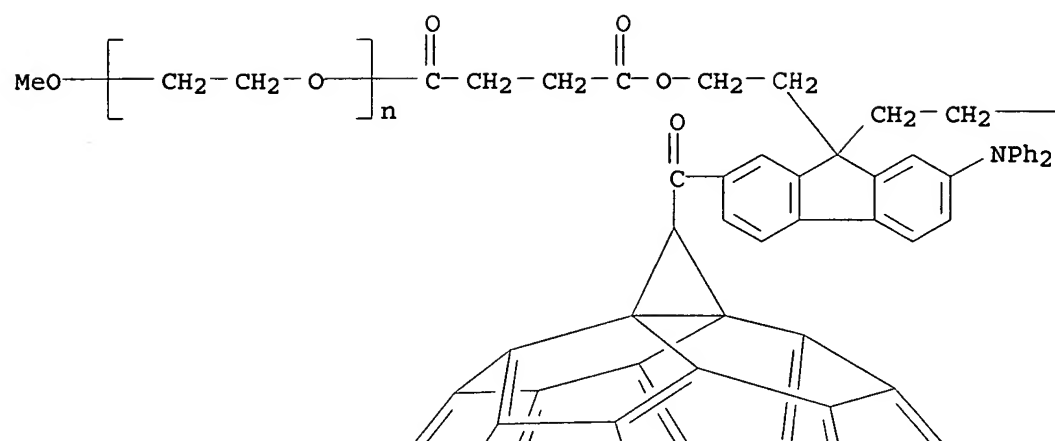
REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 11 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2005:727593 CAPLUS  
 DOCUMENT NUMBER: 144:312635  
 TITLE: Synthesis and water-solubility of pegylated diphenylaminofluorenyl monoadduct of fullerene  
 AUTHOR(S): Verma, Sarika; Padmawar, Prashant A.; Canteenwala, Taizoon; Chiang, Long Y.  
 CORPORATE SOURCE: Department of Chemistry, University of Massachusetts Lowell, Lowell, MA, 01854, USA  
 SOURCE: Polymer Preprints (American Chemical Society, Division of Polymer Chemistry) (2005), 46(2), 711-712  
 CODEN: ACPPAY; ISSN: 0032-3934  
 PUBLISHER: American Chemical Society, Division of Polymer Chemistry  
 DOCUMENT TYPE: Journal; (computer optical disk)  
 LANGUAGE: English

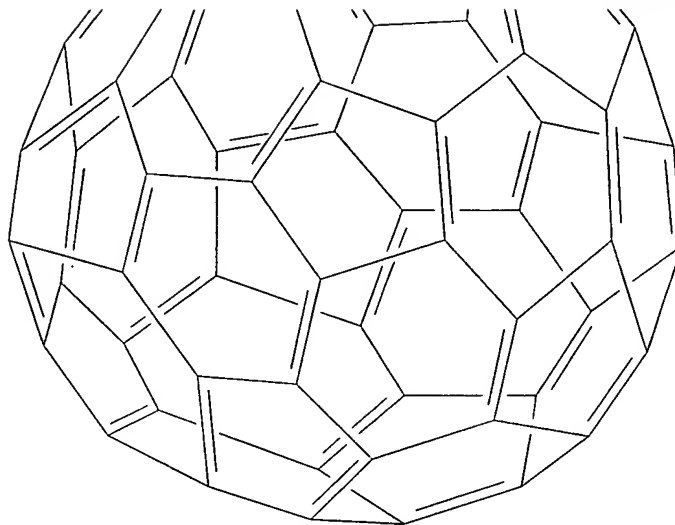
AB Recent development on two-photon absorption (2PA) excitation process may serve as a potential alternative PDT treatment procedure because of its ability to focus on a confined small treatment area of diseased tissues in a greater depth using a spectral window of 800-1100 nm in mammalian tissue. Synthesis of amphiphilic fullerene-chromophore conjugates containing polar hydrophilic functional groups, such as carboxylic acid and poly(ethylene glycol), leading to water-soluble fullerene derivs. enhances their possibility in biol. applications. Accordingly, we synthesized novel A-sp<sup>3</sup>-D type amphiphilic 1;603; fullerene-diphenylaminofluorene-poly(ethylene glycol) conjugates, namely, C60(>DPAF-PEG). Compound C60(>DPAF-PEG) consists of a fullerene cage as the electron acceptor (A) moiety and diphenylaminofluorene segment as the donor (D) moiety, showing high two-photon absorption (2PA) cross-sections in nanosecond regime. Here, we present the synthesis, characterization and mol. self-assembly of these amphiphilic two-photon active fullerenic materials in aqueous solns. Resulting amphiphilics were characterized by FT-IR, DSC, TEM, uv-visible and fluorescence spectroscopy.

IT 879872-39-6P  
 RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (synthesis and water-solubility of polyethylene glycol-diphenylaminofluorenyl monoadduct of fullerene)  
 RN 879872-39-6 CAPLUS  
 CN Poly(oxy-1,2-ethanediyl),  $\alpha,\alpha'$ -[[2-(3'H-cyclopropa[1,9] [5,6] fullerene-C60-1h-3'-ylcarbonyl)-7-(diphenylamino)-9H-fluoren-9-ylidene]bis[2,1-ethanediyl]oxy(1,4-dioxo-4,1-butanediyl)]]bis[ $\omega$ -methoxy- (9CI) (CA INDEX NAME)





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IT 649724-50-5

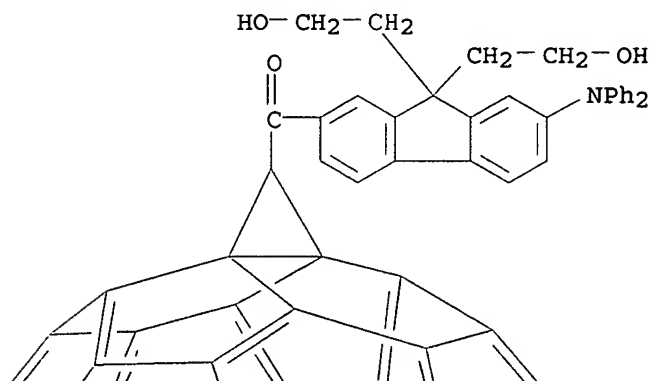
RL: RCT (Reactant); RACT (Reactant or reagent)

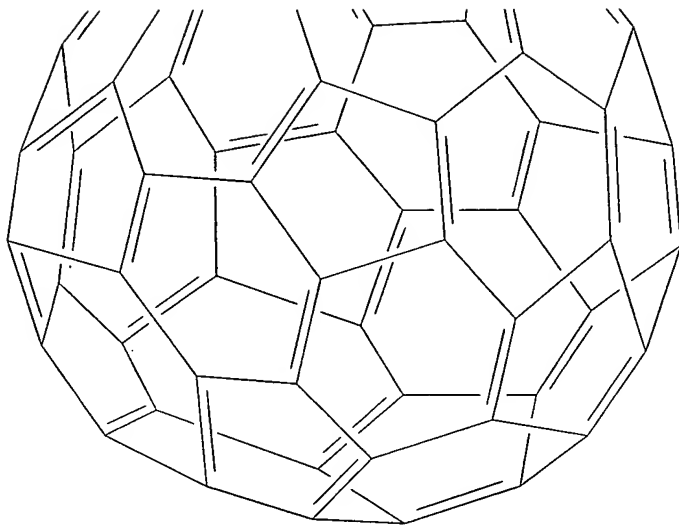
(synthesis and water-solubility of polyethylene glycol-diphenylaminofluorenyl monoadduct of fullerene)

RN 649724-50-5 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-Ih-3'-yl[7-(diphenylamino)-9,9-bis(2-hydroxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

PAGE 1-A





REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 12 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:655378 CAPLUS

DOCUMENT NUMBER: 143:306020

TITLE: Improved Synthesis of 2,2'-Dibromo-9,9'-spirobifluorene and Its 2,2'-Bisdonor-7,7'-bisacceptor-Substituted Fluorescent Derivatives

AUTHOR(S): Chiang, Chih-Long; Shu, Ching-Fong; Chen, Chin-Ti

CORPORATE SOURCE: Department of Applied Chemistry, National Chiao Tung University, Hsin-Chu, 30035, Taiwan

SOURCE: Organic Letters (2005), 7(17), 3717-3720

CODEN: ORLEF7; ISSN: 1523-7060

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 143:306020

AB Pure 2,2'-dibromo-9,9'-spirobifluorene (I) was synthesized by a method that did not involve troublesome dibromination of 9,9'-spirobifluorene or Sandmeyer reaction of 2,2'-diamino-9,9'-spirobifluorene. Starting from 4-Me<sub>3</sub>SiC<sub>6</sub>H<sub>4</sub>B(OH)<sub>2</sub>, I was prepared by Suzuki cross-coupling with 1,2-C<sub>6</sub>H<sub>4</sub>Br<sub>2</sub>, subsequent lithiation and condensation with (MeO)<sub>2</sub>CO, further bromodesilation, and finally spirocyclization by classical Clark and Gomberg method. A series of donor-acceptor orthogonally substituted 9,9'-spirobifluorene was subsequently prepared showing rich variation of fluorescence in solution and in solid state. Compound I was studied by x-ray structural anal. [monoclinic, space group P2(1)/c, a 14.5655(5), b 16.5819(5), c 7.9981(2) Å, β 93.4850(10)°, V 1928.16(10) Å<sup>3</sup>, Z 4].

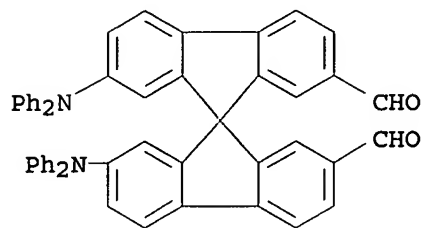
IT 864957-76-6P

RL: PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of bromospirobifluorene and bisdonor bisacceptor fluorescent derivs.)

RN 864957-76-6 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2,2'-dicarboxaldehyde, 7,7'-bis(diphenylamino)-  
(9CI) (CA INDEX NAME)

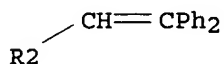
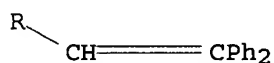
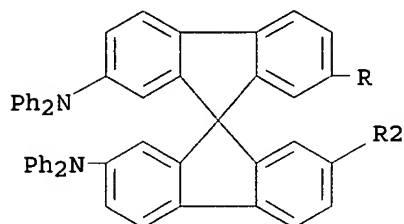


IT 724789-65-5P 864957-77-7P 864957-78-8P  
864957-79-9P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
(preparation of bromospirobifluorene and bisdonor bisacceptor fluorescent  
derivs.)

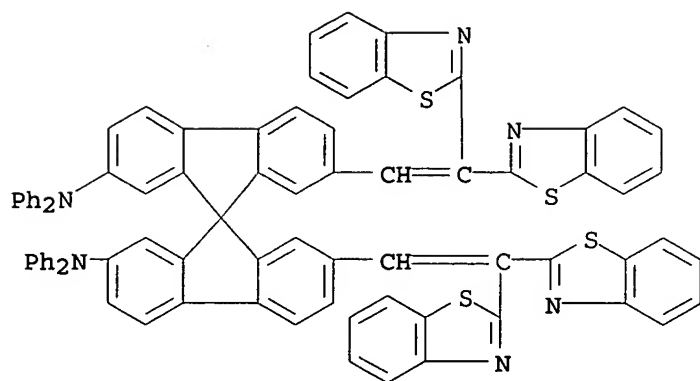
RN 724789-65-5 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2,2'-diamine, 7,7'-bis(2,2-diphenylethenyl)-  
N,N,N',N'-tetraphenyl- (9CI) (CA INDEX NAME)



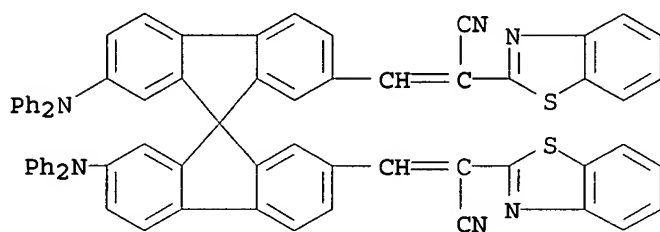
RN 864957-77-7 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2,2'-diamine, 7,7'-bis[2,2-bis(2-  
benzothiazolyl)ethenyl]-N,N,N',N'-tetraphenyl- (9CI) (CA INDEX NAME)



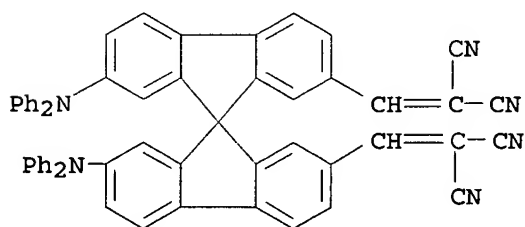
RN 864957-78-8 CAPLUS

CN 2-Benzothiazoleacetonitrile,  $\alpha,\alpha'$ -[[7,7'-bis(diphenylamino)-9,9'-spirobi[9H-fluorene]-2,2'-diyl]dimethyldiylidene]bis- (9CI) (CA INDEX NAME)



RN 864957-79-9 CAPLUS

CN Propanedinitrile, 2,2'-[[7,7'-bis(diphenylamino)-9,9'-spirobi[9H-fluorene]-2,2'-diyl]dimethyldiylidene]bis- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 38 THERE ARE 38 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 13 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:524644 CAPLUS

DOCUMENT NUMBER: 143:193797

TITLE: Synthesis of hydrophilic two-photon absorptive fullerene-diphenylaminofluorene dyads for molecular self-assembly in water

AUTHOR(S): Verma, Sarika; Hauck, Tanya; Padmawar, Prashant A.; Canteenwala, Taizoon; Chiang, Long Y.; Pritzker, Kenneth P. H.

CORPORATE SOURCE: Department of Chemistry, Institute of Nanoscience and Engineering, University of Massachusetts, Lowell, MA, 01854, USA

SOURCE: Materials Research Society Symposium Proceedings (2005), 846(Organic and Nanocomposite Optical Materials), 275-280  
CODEN: MRSPDH; ISSN: 0272-9172

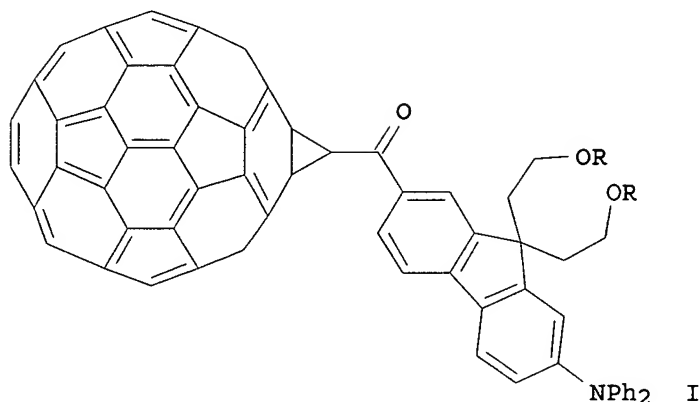
PUBLISHER: Materials Research Society

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 143:193797

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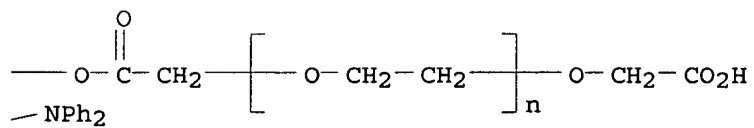
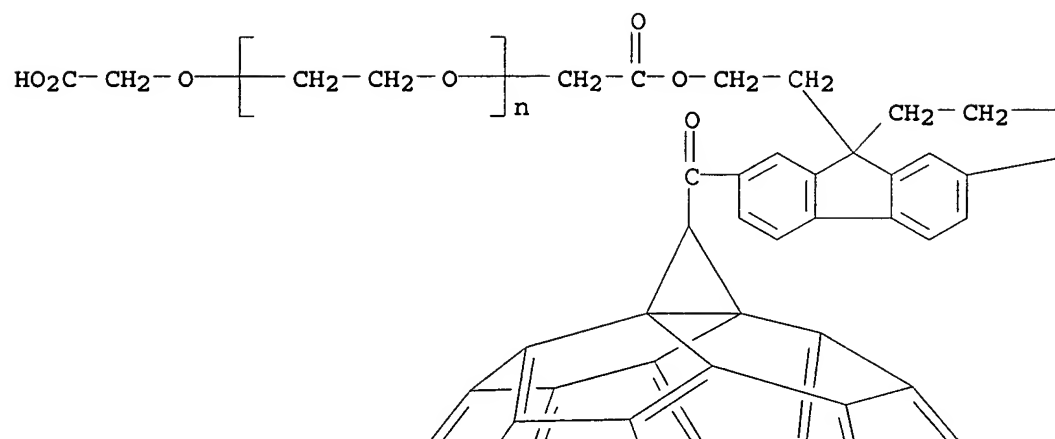


AB Diphenylaminofluorenylcarbonylfullerene I [R = HO<sub>2</sub>CCH<sub>2</sub>(OCH<sub>2</sub>CH<sub>2</sub>)<sub>n</sub>OCH<sub>2</sub>CO] is prepared by DCC-mediated esterification of poly(ethylene glycol) bis(carboxymethyl) ether with I (R = H). I [R = HO<sub>2</sub>CCH<sub>2</sub>(OCH<sub>2</sub>CH<sub>2</sub>)<sub>n</sub>OCH<sub>2</sub>CO] forms spherical aggregates 200-400 nm in diameter by self-assembly at 10 μM concentration in a chloroform/water mixture I [R = HO<sub>2</sub>CCH<sub>2</sub>(OCH<sub>2</sub>CH<sub>2</sub>)<sub>n</sub>OCH<sub>2</sub>CO] is a two-photon absorbing agent with potential for biomedical applications (no data).

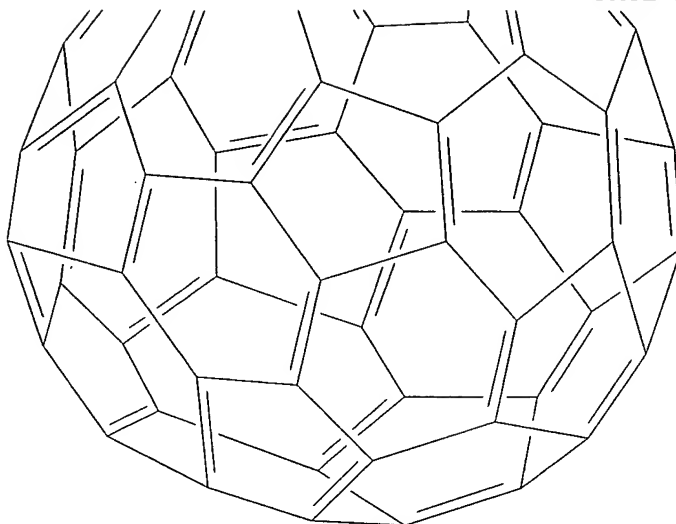
IT 851204-04-1P  
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation and self-assembly of a poly(ethylene glycol)-substituted diphenylaminofluorenylcarbonylfullerene as a hydrophilic two-photon-absorbing agent)

RN 851204-04-1 CAPLUS

CN Poly(oxy-1,2-ethanediyl), α,α'-[[2-(3'H-cyclopropa[1,9][5,6]fullerene-C<sub>60</sub>-1h-3'-ylcarbonyl)-7-(dipropylamino)-9H-fluoren-9-ylidene]bis[2,1-ethanediylxy(2-oxo-2,1-ethanediyl)]]bis[ω-(carboxymethoxy)-(9CI) (CA INDEX NAME)



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IT 649724-50-5

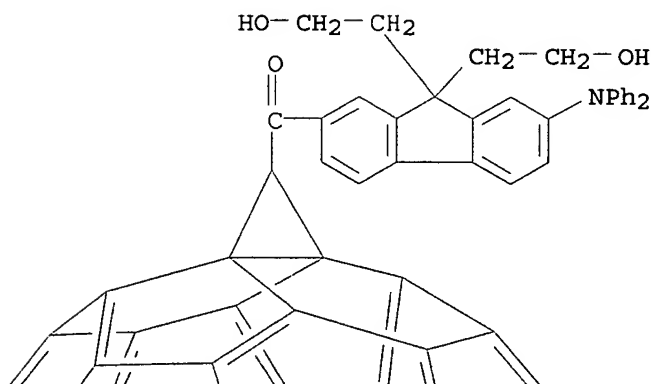
RL: RCT (Reactant); RACT (Reactant or reagent)

(preparation and self-assembly of a poly(ethylene glycol)-substituted  
diphenylaminofluorenylcarbonylfullerene as a hydrophilic  
two-photon-absorbing agent)

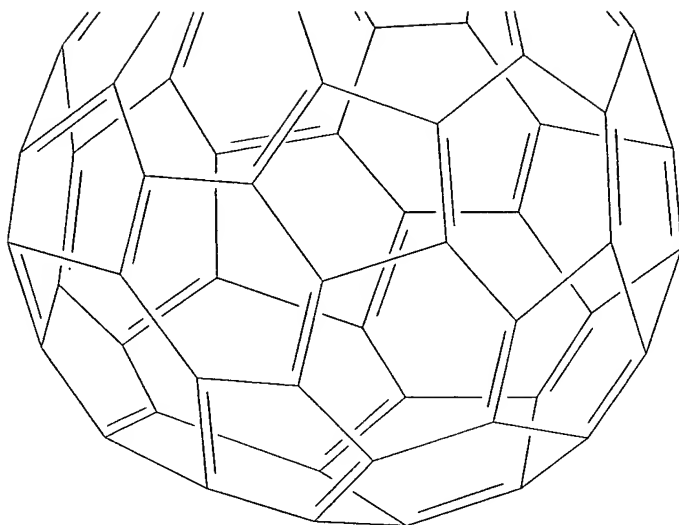
RN 649724-50-5 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-Ih-3'-yl[7-(diphenylamino)-  
9,9-bis(2-hydroxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

PAGE 1-A







REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 14 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:239950 CAPLUS

DOCUMENT NUMBER: 143:459708

TITLE: Red-emitting fluorenes as efficient emitting hosts for non-doped, organic red-light-emitting diodes

AUTHOR(S): Chiang, Chih-Long; Wu, Min-Fei; Dai, De-Chang; Wen, Yuh-Sheng; Wang, Juen-Kai; Chen, Chin-Ti

CORPORATE SOURCE: Institute of Chemistry, Academia Sinica, Taipei, 11529, Taiwan

SOURCE: Advanced Functional Materials (2005), 15(2), 231-238  
CODEN: AFMDC6; ISSN: 1616-301X

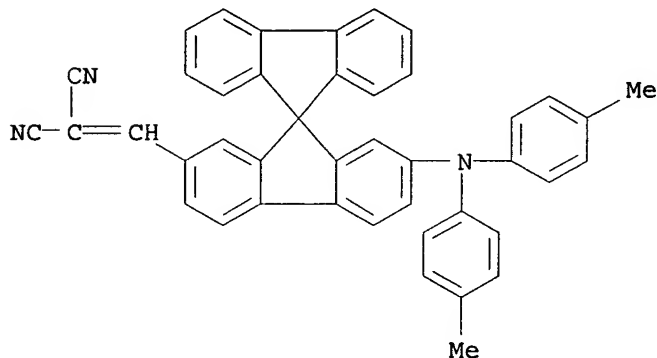
PUBLISHER: Wiley-VCH Verlag GmbH & Co. KGaA

DOCUMENT TYPE: Journal

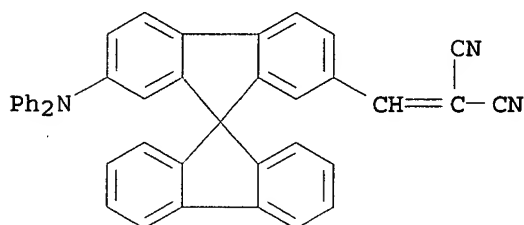
LANGUAGE: English

AB Rare red-fluorescent fluorene derivs. were designed and synthesized. The long-wavelength red fluorescence is achieved by incorporating a di(4-tolyl)amino or diphenylamino electron donor and a dicyanovinyl electron acceptor. The single-crystal x-ray structures of the di(4-tolyl)amino (pTSPDCV) and diphenylamino (PhSPDCV) compds. indicate only weak non- $\pi$  van der Waals contacts in addition to long-distance dipole-dipole interactions of the red-emitting fluorene mols. in the solid state. The aggregation of the dipolar fluorene is largely suppressed by introducing bulky 9,9-substituents (spiro-fused bifluorene) as well as a nonplanar di(4-tolyl)amino or diphenylamino group. In the solid state, these fluorene derivs. all show red fluorescence that is much brighter than with the red dopants Nile Red and DCM (4-(dicyanomethylene)-2-methyl-6-[4-(dimethylaminostyryl)-4H-pyran]). The unique photophys. properties of red-emitting fluorene derivs. differ from other known red dopants and facilitate the fabrication of nondoped red organic light-emitting diodes (OLEDs). Authentic red (CIE,  $x = 0.65$ ,  $y = 0.35$ ) electroluminescence with a brightness of  $>12000$  cd  $m^{-2}$  (greater than 600 cd  $m^{-2}$  at 20 mA  $cm^{-2}$ ) and a remarkable external quantum efficiency  $\leq 3.6\%$  were observed for

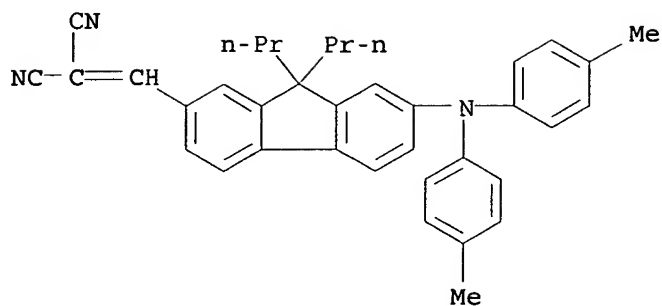
the red-emitting OLEDs with pTSPDCV or PhSPDCV as the sole emitting host.  
 IT 869299-85-4P 869299-86-5P  
 RL: DEV (Device component use); PRP (Properties); SPN (Synthetic preparation); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)  
 (crystallog. and red fluorescence; red-emitting fluorenes as efficient emitting hosts for non-doped, organic red-light-emitting diodes)  
 RN 869299-85-4 CAPLUS  
 CN Propanedinitrile, [[7-[bis(4-methylphenyl)amino]-9,9'-spirobi[9H-fluoren]-2-yl]methylene]- (9CI) (CA INDEX NAME)



RN 869299-86-5 CAPLUS  
 CN Propanedinitrile, [[7-(diphenylamino)-9,9'-spirobi[9H-fluoren]-2-yl]methylene]- (9CI) (CA INDEX NAME)



IT 869299-84-3P  
 RL: DEV (Device component use); PRP (Properties); SPN (Synthetic preparation); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)  
 (red fluorescence; red-emitting fluorenes as efficient emitting hosts for non-doped, organic red-light-emitting diodes)  
 RN 869299-84-3 CAPLUS  
 CN Propanedinitrile, [[7-[bis(4-methylphenyl)amino]-9,9-dipropyl-9H-fluoren-2-yl]methylene]- (9CI) (CA INDEX NAME)



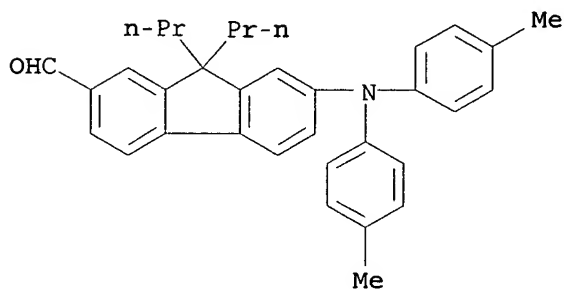
IT 869299-87-6P 869299-88-7P 869299-89-8P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(red-emitting fluorenes as efficient emitting hosts for non-doped, organic red-light-emitting diodes)

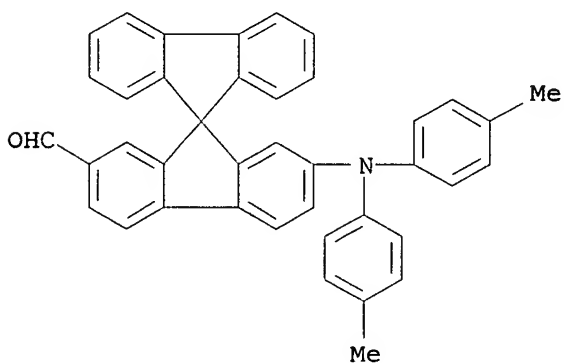
RN 869299-87-6 CAPLUS

CN 9H-Fluorene-2-carboxaldehyde, 7-[bis(4-methylphenyl)amino]-9,9-dipropyl- (9CI) (CA INDEX NAME)



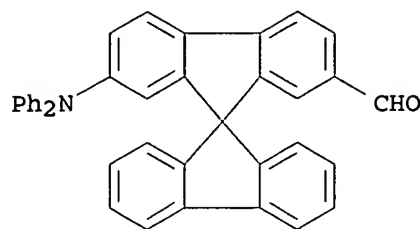
RN 869299-88-7 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2-carboxaldehyde, 7-[bis(4-methylphenyl)amino]- (9CI) (CA INDEX NAME)



RN 869299-89-8 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2-carboxaldehyde, 7-(diphenylamino)- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 47 THERE ARE 47 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 15 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:224493 CAPLUS

DOCUMENT NUMBER: 142:454178

TITLE: Self-Assembled Photoresponsive Amphiphilic Diphenylaminofluorene-C60 Conjugate Vesicles in Aqueous Solution

AUTHOR(S): Verma, Sarika; Hauck, Tanya; El-Khouly, Mohamed E.; Padmawar, Prashant A.; Canteenwala, Taizoon; Pritzker, Kenneth; Ito, Osamu; Chiang, Long Y.

CORPORATE SOURCE: Department of Chemistry, Institute of Nanoscience and Engineering, University of Massachusetts, Lowell, MA, 01854, USA

SOURCE: Langmuir (2005), 21(8), 3267-3272

CODEN: LANGD5; ISSN: 0743-7463

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

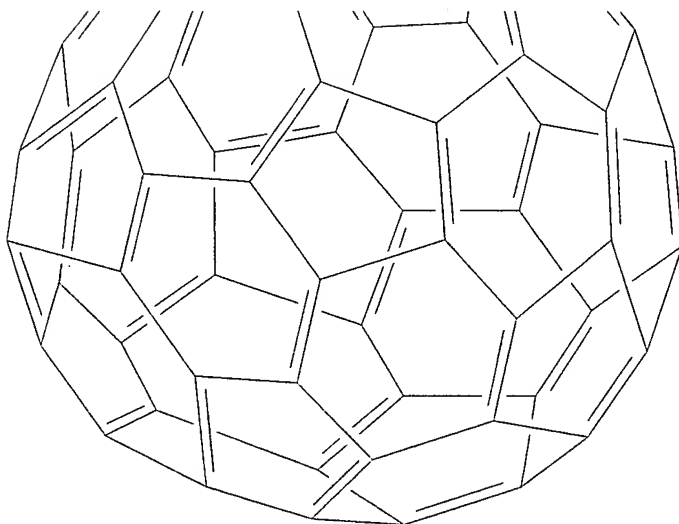
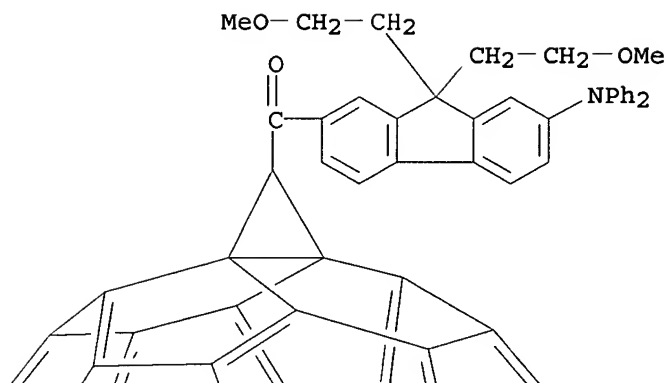
AB Water-soluble oligo(ethylene glycolated) derivs. of two-photon absorptive diphenylaminofluorencarbonyl-methano[60]fullerene, denoted as C60(> DPAF-EG6), were synthesized with their mol. self-assembly characteristics in H2O studied. The formation of nano- to submicron-sized spherical hollow vesicles with a shell width of 15-20 nm was observed by transmission electron microscopy (TEM) micrographs. This shell width fits approx. with the length of a disordered bilayer-like mol. packing of C60(> DPAF-EG6), arising from strong intermol. hydrophobic interactions of fullerene cages. Photoinduced intramol. charge separation followed by charge recombination on the nanosecond time scale, from the DPAF moiety to the C60 cage in the vesicle structure, was detected via transient spectroscopic measurements. Electronic supplementary information (ESI) is available at <http://pubs.acs.org> and contains NMR spectra and IR spectra and TGA profiles of DPAF-EG6 and its precursor compds.

IT 649724-49-2

RL: RCT (Reactant); RACT (Reactant or reagent)  
(demethylation with BBr3 in presence of NaI and 15-crown-5)

RN 649724-49-2 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

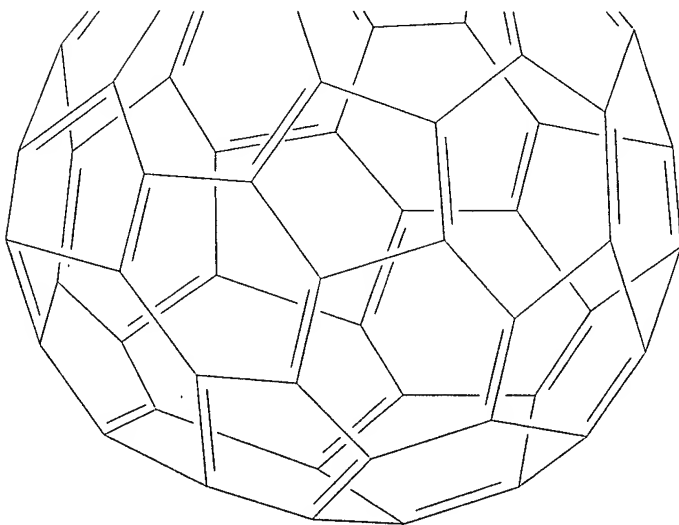
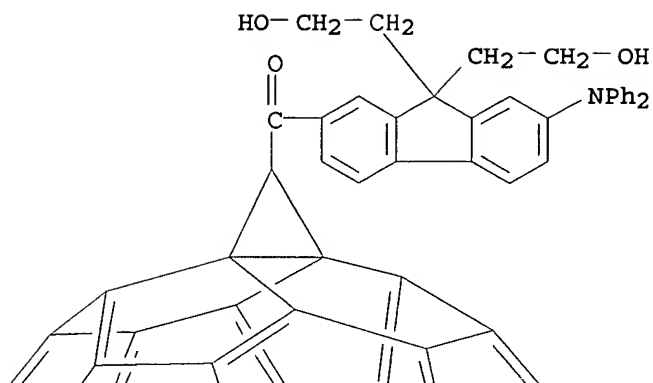


IT 649724-50-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(esterification with oligo(ethylene glycol)-bis(carboxymethyl)ether)

RN 649724-50-5 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-bis(2-hydroxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



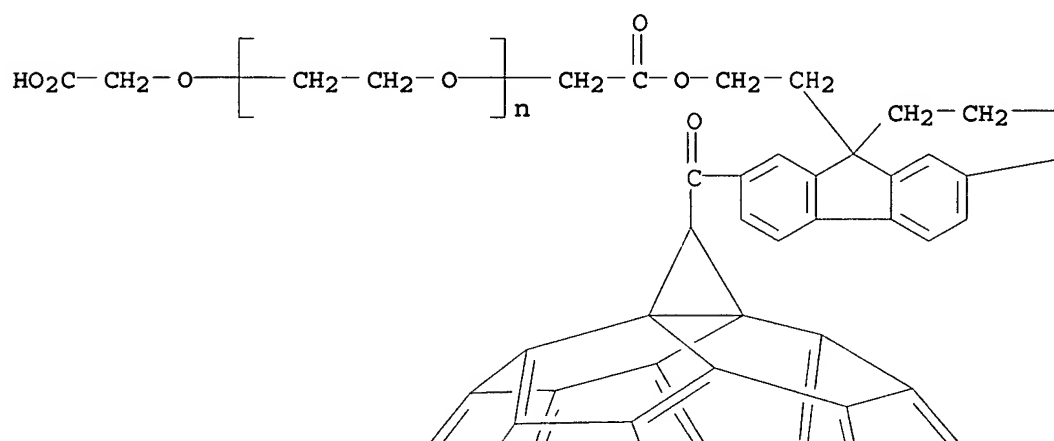
IT 851204-04-1P

RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

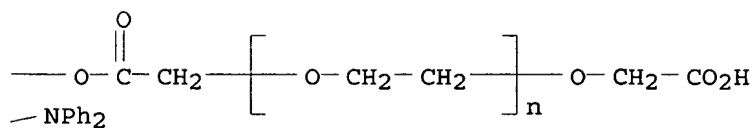
(photophys. of oligo(ethylene glycolated) derivs. of two-photon absorbing diphenylaminofluorenylcarbonyl-methano[60]fullerene in vesicle

structure)  
 RN 851204-04-1 CAPLUS  
 CN Poly(oxy-1,2-ethanediyl),  $\alpha, \alpha'$ -[[2-(3'H-cyclopropa[1,9][5,6]fullerene-C60-Ih-3'-ylcarbonyl)-7-(dipropylamino)-9H-fluoren-9-ylidene]bis[2,1-ethanediyl]oxy(2-oxo-2,1-ethanediyl)]]bis[ $\omega$ -(carboxymethoxy)-(9CI) (CA INDEX NAME)

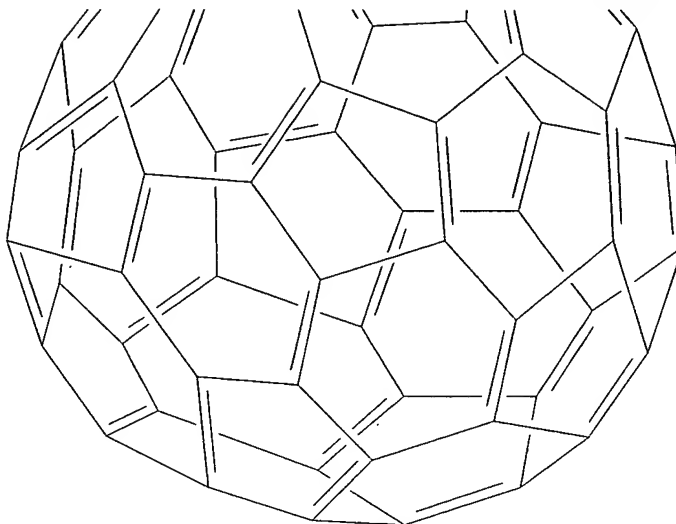
PAGE 1-A



PAGE 1-B



PAGE 2-A



REFERENCE COUNT: 27 THERE ARE 27 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 16 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:127098 CAPLUS

DOCUMENT NUMBER: 143:468290

TITLE: One- and Two-Photon Fluorescence Anisotropy of Selected Fluorene Derivatives

AUTHOR(S): Belfield, K. D.; Bondar, M. V.; Hales, J. M.; Morales, A. R.; Przhonska, O. V.; Schafer, K. J.

CORPORATE SOURCE: Department of Chemistry, University of Central Florida, Orlando, FL, USA

SOURCE: Journal of Fluorescence (2005), 15(1), 3-11

CODEN: JOFLEN; ISSN: 1053-0509

PUBLISHER: Springer

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The steady-state excitation anisotropy spectra of fluorene derivs. were measured in viscous solvents, under the one- and two-photon excitation, over a broad spectral range (UV-Visible). The orientation of their absorption transition moments for the first,  $S_0 \rightarrow S_1$ , and second,  $S_0 \rightarrow S_2$ , excited states were determined. It was shown exptl. that a decrease in the angle between  $S_0 \rightarrow S_1$  and  $S_0 \rightarrow S_2$  transitions corresponded to an increased value of two-photon absorption (2PA) cross section for these mols. Two-photon excitation anisotropy was nearly constant over the spectral region investigated (in contrast to one-photon excitation anisotropy spectra) and can be roughly explained by a simple model of 2PA based on the single intermediate state approximation. For comparison, the same trend in two-photon excitation anisotropy was observed for Rhodamine B in glycerol.

IT 252000-05-8

RL: PRP (Properties)

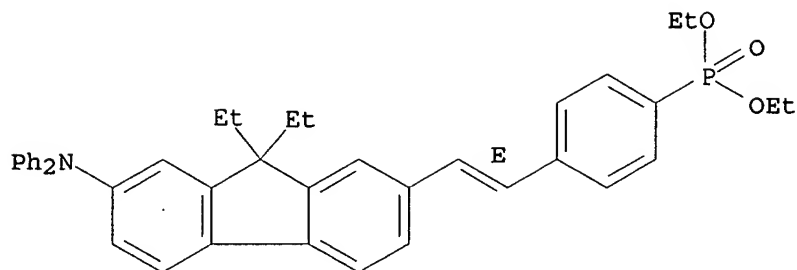
(one- and two-photon fluorescence anisotropy of selected fluorene derivs.)

RN 252000-05-8 CAPLUS

CN Phosphonic acid, [4-[(1E)-2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)

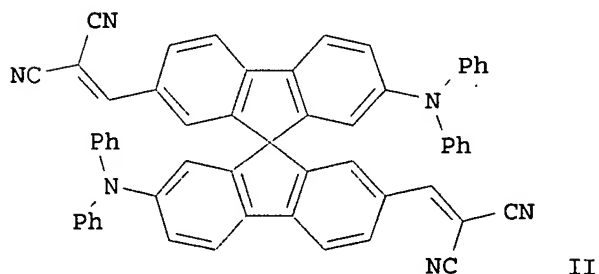
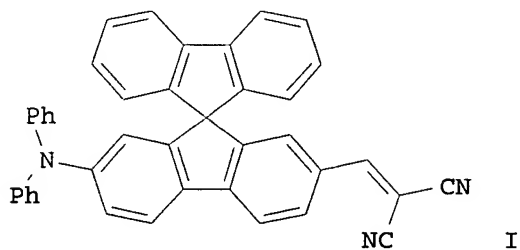


Double bond geometry as shown.



REFERENCE COUNT: 31 THERE ARE 31 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 17 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2005:114653 CAPLUS  
 DOCUMENT NUMBER: 144:36070  
 TITLE: Red fluorenes as the efficient host emitter for non-doped red organic light-emitting diodes  
 AUTHOR(S): Chiang, Chih-Long; Wu, Min-Fei; Shu, Ching-Fong; Chen, Chin-Ti  
 CORPORATE SOURCE: Department of the Applied Chemistry, National Chiao Tung Univ., Hsinchu, 30035, Taiwan  
 SOURCE: Proceedings of SPIE-The International Society for Optical Engineering (2005), 5632(Light-Emitting Diode Materials and Devices), 80-87  
 CODEN: PSISDG; ISSN: 0277-786X  
 PUBLISHER: SPIE-The International Society for Optical Engineering  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 GI



AB Crystal red fluorophores based on donor-acceptor substituted spirofluorene, i.e., I show strong fluorescence in solution ( $\Phi_f$  apprx. 70%) as well as in solid state ( $\Phi_f > 30\%$ ). Non-doped red OLEDs fabricated with I exhibit authentic red (CIE,  $x = 0.65$ ,  $y_r = 0.35$ ) electroluminescence with brightness over 12,000 cd m<sup>-2</sup> (or > 600 cd m<sup>-2</sup> at 20 mA cm<sup>-2</sup>) and remarkable external quantum efficiency as high as 3.6%. On the other hand, the bis-substituted derivs. of spirofluorene II show relatively weak fluorescence both in solution ( $\Phi_f < 20\%$ ) and in solid state ( $\Phi_f < 10\%$ ). Although saturated red electroluminescence (CIE,  $x = 0.65$ ,  $y_r = 0.34$ ) is also observed, non-doped red OLED containing II performs

much worse than I OLEDs. Both PhSPDCV and BisPhSPDCV are not amorphous forming loosely packed crystal materials in solid state with no intimate  $\pi$ - $\pi$  interaction.

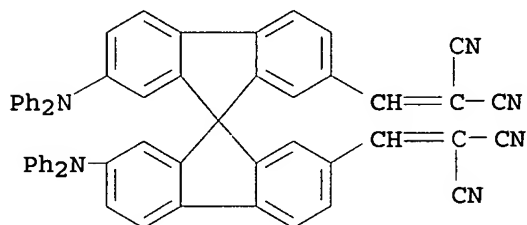
IT 864957-79-9P 869299-86-5P

RL: DEV (Device component use); PRP (Properties); SPN (Synthetic preparation); PREP (Preparation); USES (Uses)

(preparation of red fluorenes as efficient host emitter for non-doped red organic light-emitting diodes)

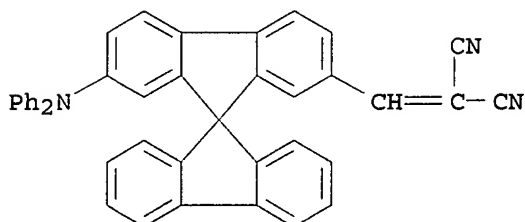
RN 864957-79-9 CAPLUS

CN Propanedinitrile, 2,2'-[[7,7'-bis(diphenylamino)-9,9'-spirobi[9H-fluorene]-2,2'-diyl]dimethylidyne]bis- (9CI) (CA INDEX NAME)



RN 869299-86-5 CAPLUS

CN Propanedinitrile, [[7-(diphenylamino)-9,9'-spirobi[9H-fluorene]-2-yl]methylene]- (9CI) (CA INDEX NAME)



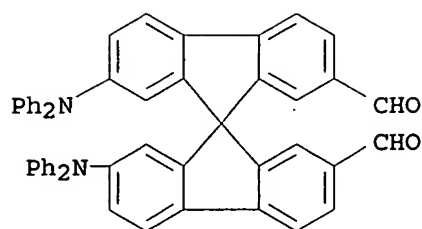
IT 864957-76-6P 869299-89-8P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

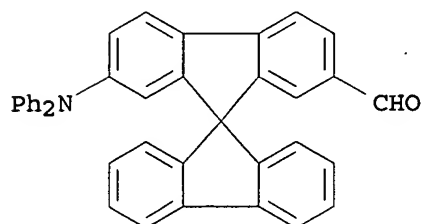
(preparation of red fluorenes as efficient host emitter for non-doped red organic light-emitting diodes)

RN 864957-76-6 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2,2'-dicarboxaldehyde, 7,7'-bis(diphenylamino)- (9CI) (CA INDEX NAME)



RN 869299-89-8 CAPLUS  
 CN 9,9'-Spirobi[9H-fluorene]-2-carboxaldehyde, 7-(diphenylamino)- (9CI) (CA  
 INDEX NAME)



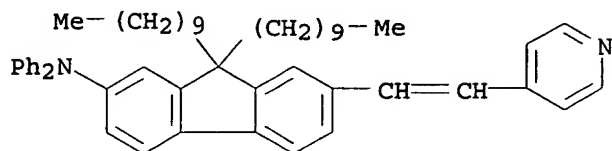
REFERENCE COUNT: 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 18 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2005:45072 CAPLUS  
 DOCUMENT NUMBER: 142:305594  
 TITLE: TDDFT Study of One- and Two-Photon Absorption  
 Properties: Donor- $\pi$ -Acceptor Chromophores  
 AUTHOR(S): Day, Paul N.; Nguyen, Kiet A.; Pachter, Ruth  
 CORPORATE SOURCE: Materials and Manufacturing Directorate, Air Force  
 Research Laboratory, Wright Patterson Air Force Base,  
 OH, 45433, USA  
 SOURCE: Journal of Physical Chemistry B (2005), 109(5),  
 1803-1814  
 CODEN: JPCBFK; ISSN: 1520-6106  
 PUBLISHER: American Chemical Society  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English

AB We report a comprehensive time-dependent d. functional theory (TDDFT)  
 study of one-photon and two-photon absorption (OPA and TPA, resp.) spectra  
 for donor- $\pi$ -acceptor mols. The calculated excitation energies were  
 generally shown to be in good agreement with experiment, particularly when  
 compared to results from measurements carried out in a nonpolar solvent,  
 although the oscillator strengths were overestimated in some cases.  
 Calculated TPA cross sections applying the two-state approximation were shown  
 to be highly dependent on the form of the line-shape function used. Although a  
 good agreement with exptl. TPA spectra was generally observed, the wide range  
 in the exptl. measured values and lack of systematic exptl. data on  
 solvent effects limited a detailed comparison as yet.

IT 191667-13-7 252000-05-8  
 RL: PRP (Properties)  
 (TDDFT study of one- and two-photon absorption of donor- $\pi$ -acceptor  
 chromophores)  
 RN 191667-13-7 CAPLUS

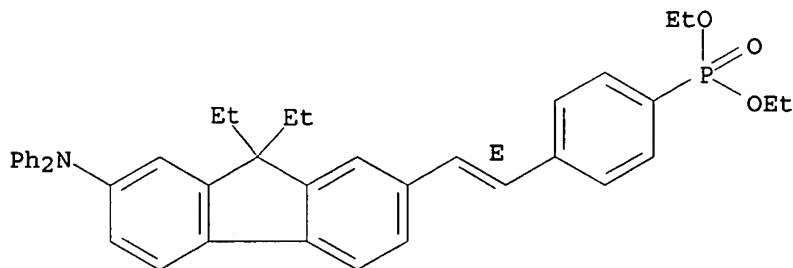
CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



RN 252000-05-8 CAPLUS

CN Phosphonic acid, [4-[(1E)-2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



REFERENCE COUNT: 104 THERE ARE 104 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE REFORMAT

L6 ANSWER 19 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:3217 CAPLUS

DOCUMENT NUMBER: 143:294805

TITLE: Substituent effect on two-photon absorption properties of fluorene derivatives

AUTHOR(S): Kawamata, Jun; Akiba, Masaharu; Inagaki, Yoshio; Tani, Takeharu; Harada, Akinori

CORPORATE SOURCE: Department of Chemistry, Faculty of Science, Yamaguchi University, Yamaguchi, 753-8512, Japan

SOURCE: Journal of Nonlinear Optical Physics & Materials (2004), 13(3 & 4), 475-479

CODEN: JNOMFV; ISSN: 0218-8635

PUBLISHER: World Scientific Publishing Co. Pte. Ltd.

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Two novel fluorene derivs. having cationic substituents were synthesized. Two-photon absorption (TPA) properties of the derivs. were evaluated using the fluorescence-based technique with a femtosecond pulse emitted from a Ti:sapphire laser. Maximum TPA cross sections of the compds. were estimated from the TPA spectra. Based on the results, the relation between the TPA cross sections and the substituent effects of the derivs. are discussed.

IT 197969-56-5 437713-15-0

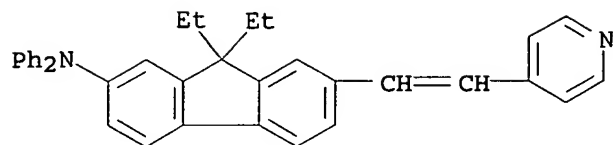
RL: PRP (Properties)

(substituent effect on two-photon absorption properties of fluorene derivs.)

RN 197969-56-5 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-

(9CI) (CA INDEX NAME)



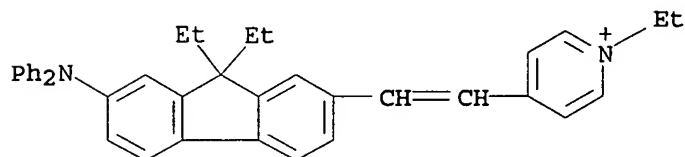
RN 437713-15-0 CAPLUS

CN Pyridinium, 4-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]-1-ethyl-, perchlorate (9CI) (CA INDEX NAME)

CM 1

CRN 437713-14-9

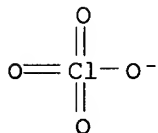
CMF C38 H37 N2



CM 2

CRN 14797-73-0

CMF Cl O4



REFERENCE COUNT: 13 THERE ARE 13 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 20 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:1149754 CAPLUS

DOCUMENT NUMBER: 143:237948

TITLE: Nonlinear transmission and excited-state absorption in fluorene derivatives

AUTHOR(S): Belfield, Kevin D.; Bondar, Mykhailo V.; Hernandez, Florencio E.; Morales, Alma R.; Przhonska, Olga V.; Schafer, Katherine J.

CORPORATE SOURCE: Department of Chemistry and College of Optics and Lasers and the Florida Photonics Center of Excellence, University of Central Florida, Orlando, FL, 32816-2366, USA

SOURCE: Applied Optics (2004), 43(34), 6339-6343  
CODEN: APOPAI; ISSN: 0003-6935

PUBLISHER: Optical Society of America

DOCUMENT TYPE: Journal  
 LANGUAGE: English

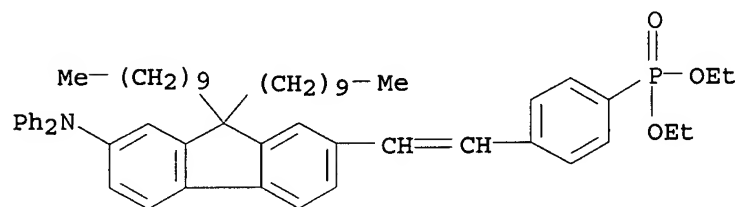
AB The nonlinear transmission and the excited-state absorption spectra of 3 fluorene derivs. exhibiting large 2-photon absorptivity were measured by the 3rd harmonic of a picosecond Nd:YAG laser. We analyzed their capability for exhibiting stimulated emission in polar solvents and found that asym. fluorene compds. with a diphenylamino substituent exhibited large Stokes shifts (.apprx.8000 cm<sup>-1</sup>), high quantum yields (.apprx.0.9-1.0), and no optical gain over their entire fluorescence spectral region. In contrast, a sym. fluorene derivative with vinylphenylbenzothiazole substituents in positions 2 and 7 underwent lasing under 1-photon excitation by use of picosecond pulsed irradiation

IT 507271-13-8

RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); PROC (Process)  
 (nonlinear transmission and excited-state absorption in fluorene derivs.)

RN 507271-13-8 CAPLUS

CN Phosphonic acid, [4-[2-[9,9-didecyl-7-(diphenylamino)-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)



REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 21 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:962266 CAPLUS

DOCUMENT NUMBER: 142:143893

TITLE: Synthesis and Photophysical Properties of C60-Diphenylaminofluorene Dyad and Multiads

AUTHOR(S): Padmawar, Prashant A.; Canteenwala, Taizoon; Verma, Sarika; Tan, Loon-Seng; Chiang, Long Y.

CORPORATE SOURCE: Department of Chemistry, Institute of Nanoscience and Engineering Technology, University of Massachusetts Lowell, Lowell, MA, USA

SOURCE: Journal of Macromolecular Science, Pure and Applied Chemistry (2004), A41(12), 1387-1400  
 CODEN: JSPCE6; ISSN: 1060-1325

PUBLISHER: Marcel Dekker, Inc.

DOCUMENT TYPE: Journal

LANGUAGE: English

AB One mono-adduct and three novel multi-adduct analogs of C60-diphenylaminofluorene C60(>DPAF)<sub>n</sub> derivs. were synthesized and characterized by various spectroscopic methods. Optical absorption of these samples indicated a systematic increase in relative intensity of the fluorene band centered at 415 nm as the number of DPAF addend increases. Fluorescence of DPAF chromophore in all C60(>DPAF)<sub>n</sub> derivs. 4 (n = 1), 5 (n = 2), 6 (n = 3), and 7 (n = 4) in o-dichlorobenzene and chloroform was found to be efficiently quenched as the direct covalent bond attachment of DPAF moieties to the fullerene cage facilitates efficient intramol. electron or energy transfer processes.

IT 486998-58-7P

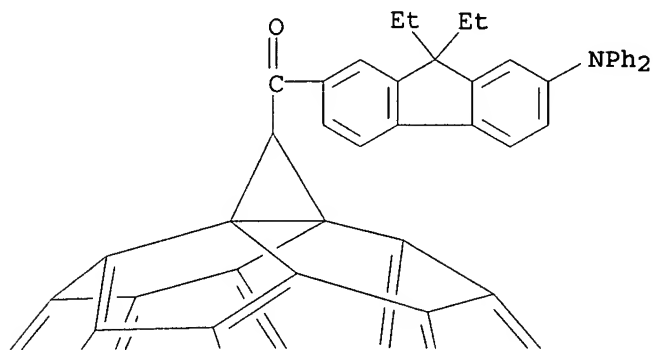
RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

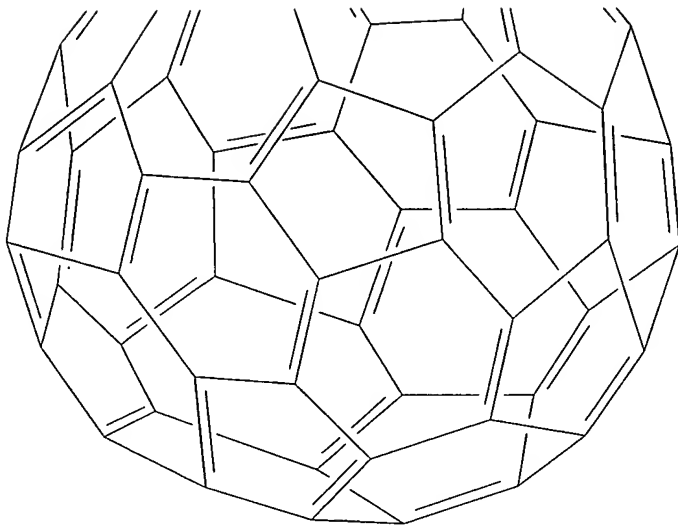
(dyad; synthesis and material characterization and photophys. properties of fullerene-diphenylaminofluorene dyad and its multi-adduct analogs)

RN 486998-58-7 CAPLUS

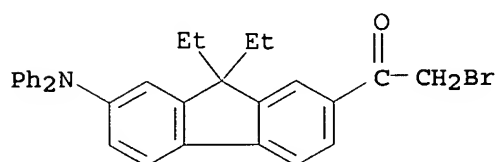
CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-Ih-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

PAGE 1-A

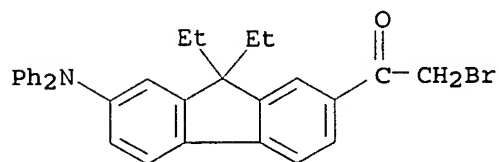




IT 486998-57-6DP, bis-, tris-, or tetra- cyclopropylated with C60  
 RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)  
 (synthesis and material characterization and photophys. properties of fullerene-diphenylaminofluorene dyad and its multi-adduct analogs)  
 RN 486998-57-6 CAPLUS  
 CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI)  
 (CA INDEX NAME)



IT 486998-57-6P, 7-Bromoacetyl-9,9-Diethyl-2-diphenylaminofluorene  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (synthesis of fullerene-diphenylaminofluorene dyad)  
 RN 486998-57-6 CAPLUS  
 CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI)  
 (CA INDEX NAME)



REFERENCE COUNT: 31 THERE ARE 31 CITED REFERENCES AVAILABLE FOR THIS

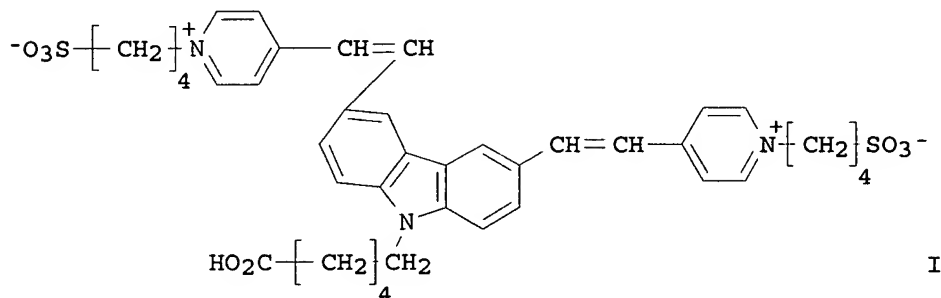


RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 22 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2004:651328 CAPLUS  
 DOCUMENT NUMBER: 141:170427  
 TITLE: Two-photon absorbing compounds and method for  
 fluorescent labeling with the compounds  
 INVENTOR(S): Inagaki, Yoshio; Takizawa, Hiroo; Akiba, Masaharu  
 PATENT ASSIGNEE(S): Fuji Photo Film Co., Ltd., Japan  
 SOURCE: Jpn. Kokai Tokkyo Koho, 14 pp.  
 CODEN: JKXXAF  
 DOCUMENT TYPE: Patent  
 LANGUAGE: Japanese  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2004224746	A2	20040812	JP 2003-15683	20030124
PRIORITY APPLN. INFO.:			JP 2003-15683	20030124
OTHER SOURCE(S):	MARPAT	141:170427		

GI



AB Substances are labeled by reacting with two-photon absorbing compds. represented by G(L)nA (G = reactive group forming a covalent bond with the substances; L = linking group; A = organic group derived from HA having two-photon absorption cross section of  $\geq 1000$  GM; n = 0, 1). Fluorometry using the two-photon absorbing compds. shows large difference between excitation wavelength and fluorescence wavelength, good three-dimensional spatial resolution, and high sensitivity. Two-photon fluorescence ratio of some compds. were given. Thus, an aqueous gelatin solution

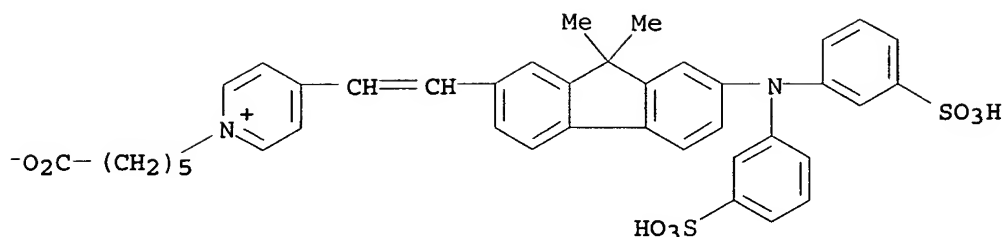
was treated with a carbazole derivative having 1 CO<sub>2</sub>H and 2 SO<sub>3</sub><sup>-</sup> groups (represented by I) and irradiated with 780-nm laser in the dark to generate bluish-green fluorescence only near the focal point.

IT 733805-44-2

RL: ARG (Analytical reagent use); PRP (Properties); ANST (Analytical study); USES (Uses)  
 (two-photon absorbing compds. and method for two-photon fluorescent labeling with the compds.)

RN 733805-44-2 CAPLUS

CN Pyridinium, 4-[2-[7-[bis(3-sulfophenyl)amino]-9,9-dimethyl-9H-fluoren-2-yl]ethenyl]-1-(5-carboxypentyl)-, inner salt, dipotassium salt (9CI) (CA INDEX NAME)



● 2 K

L6 ANSWER 23 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:646022 CAPLUS

DOCUMENT NUMBER: 142:214531

TITLE: Spectroscopic studies of the interaction between DNA and fluorene dyes

AUTHOR(S): Przonska, O. V.; Bondar, M. V.; Dovbeshko, G. I.; Repnytska, O. P.; Lukashenko, V. I.; Tryndiak, V. P.; Todor, I. N.; Belfield, K. D.

CORPORATE SOURCE: Institute of Physics, Nat. Acad. Sci. of Ukraine, Kiev, 02028, Ukraine

SOURCE: Ukrains'kii Fizichnii Zhurnal (2004), 49(7), 636-647  
CODEN: UFZHFY; ISSN: 0372-400X

PUBLISHER: Natsional'na Akademiya Nauk Ukraini, Viddilennya Fiziki i Astronomii

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Making use of electron spectroscopy (absorption and fluorescence) and IR vibrational spectroscopy methods, the interaction of the DNA taken from the normal and tumor (the yield strain of the Guerin carcinoma) tissues of rats of the Vistar line with two types of fluorene dyes, namely, (7-benzothiazol-2-yl-9,9-didecylfluoren-2-yl) diphenylamine (FD1) and {4-[2-(7-diphenylamino-9,9-diethylfluoren-2-yl)vinyl]phenyl} of the di-Et ester of phosphoric acid (FD2), has been investigated. The addition of the dyes to the tumor DNA in the in vivo experiment has been shown to disrupt the DNA secondary structure and to induce the transition into a partially unpaired form. In the IR experiment, different linkages of both dyes with the sugar-phosphate DNA frame have been detected, but it did not result in the features of the fluorescence spectra in the FD1 case. The fluorescence spectra of the FD2 interacting with the reference DNA show two bands, in contrast to what is measured at the interaction with the tumor DNA. The appearance of the short-wave band may be stipulated by an essential reconstruction in the part of the dye mols. due to their interaction with DNA ones. Exptl. results evidence for that the dyes link more efficiently to the reference DNA than to the tumor one, and that the FD2 dye with a strong acceptor group reveals an enhanced activity during the interaction. The in vivo experiment shows that addnl. sites appear in the dyes for their linkage with DNA.

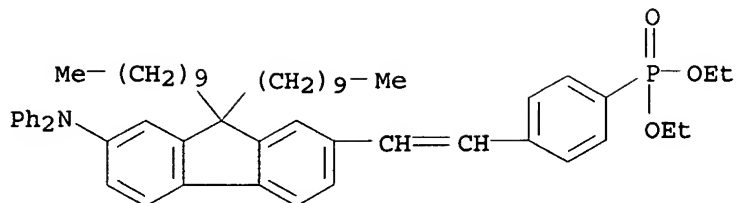
IT 507271-13-8

RL: PRP (Properties)

(spectroscopic studies of interaction between DNA and fluorene dyes)

RN 507271-13-8 CAPLUS

CN Phosphonic acid, [4-[2-[9,9-didecyl-7-(diphenylamino)-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)



L6 ANSWER 24 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:587037 CAPLUS

DOCUMENT NUMBER: 141:131068

TITLE: Electroluminescent compositions, and their organic electroluminescent devices emitting light from green to yellow

INVENTOR(S): Onikubo, Shunichi; Yauchi, Hiroyuki; Yagi, Tamao; Kaneko, Tetsuya; Tanaka, Hiroaki; Takada, Yasuyuki

PATENT ASSIGNEE(S): Toyo Ink Mfg. Co., Ltd., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 67 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent

LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2004206893	A2	20040722	JP 2002-371262	20021224

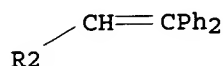
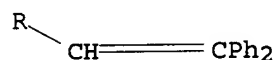
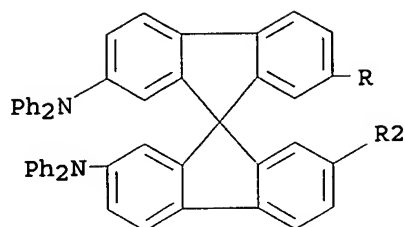
PRIORITY APPLN. INFO.: JP 2002-371262 20021224

AB The compns. contain (A) compds. having peaks at 475-600 nm in fluorescent spectra of their solid films and (B) compds. showing the sum of areas (intensities)  $\leq 20\%$  at  $\leq 500$  nm and  $\geq 600$  nm, or at  $\geq 500$  nm based on total areas (intensities) at 400-800 nm in fluorescent spectrum of solid films comprising A and 5% B. Organic electroluminescent devices having emitter layers containing the compns. containing 1:0.1 perylene derivative and diketopyrrolopyrrole derivative showed high luminescence intensity and good durability in repeated use.

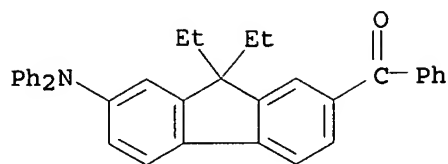
IT 724789-65-5  
 RL: DEV (Device component use); TEM (Technical or engineered material use); USES (Uses)  
 (host; electroluminescent compns. for organic electroluminescent devices showing high luminescence intensity and durability in repeated use)

RN 724789-65-5 CAPLUS

CN 9,9'-Spirobi[9H-fluorene]-2,2'-diamine, 7,7'-bis(2,2-diphenylethenyl)-N,N,N',N'-tetraphenyl- (9CI) (CA INDEX NAME)



L6 ANSWER 25 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2004:577508 CAPLUS  
 DOCUMENT NUMBER: 141:267426  
 TITLE: Few-states models for three-photon absorption  
 AUTHOR(S): Cronstrand, Peter; Norman, Patrick; Luo, Yi; Agren, Hans  
 CORPORATE SOURCE: Theoretical Chemistry, SCFAB, Royal Institute of Technology, Stockholm, SE-106 91, Swed.  
 SOURCE: Journal of Chemical Physics (2004), 121(5), 2020-2029  
 CODEN: JCPSA6; ISSN: 0021-9606  
 PUBLISHER: American Institute of Physics  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 AB Few-states models are derived for the calcn. of three-photon absorption matrix elements. Together with earlier derived few-states models for two-photon absorption, the models are evaluated against results from response theory calcns. that provide the full sum-over-states values. It is demonstrated that not even for systems with charge-transfer character, where few-states models for two-photon absorption are in excellent agreement with response theory, do the models provide a quant. correct description for three-photon absorption. The convergence behavior, merits, and shortcomings of the models are elucidated in some detail. The role of various characteristics of the electronic structure, such as symmetry, charge transfer, and conjugation-important for the formation of a large three-photon cross section-is analyzed. As for two-photon absorption cross sections, it is essential to consider generalized few-states models also for three-photon absorption, i.e., to account for dipolar directions and laser beam polarization. Despite their poor quant. performance, it is argued that few-states models at times can be useful for interpretation purposes when applied to three-photon absorption.  
 IT 349123-45-1  
 RL: PRP (Properties)  
 (few-states models for three-photon absorption in)  
 RN 349123-45-1 CAPLUS  
 CN Methanone, [7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]phenyl- (9CI)  
 (CA INDEX NAME)

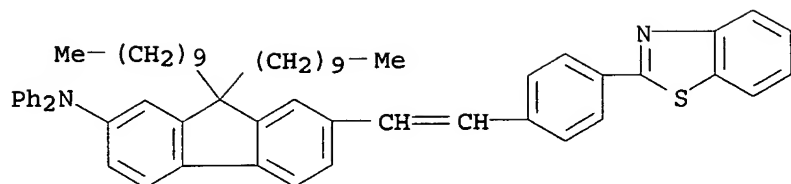


REFERENCE COUNT: 28 THERE ARE 28 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 26 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2004:543643 CAPLUS  
 DOCUMENT NUMBER: 141:313812  
 TITLE: Synthesis, Characterization, and Optical Properties of New Two-Photon-Absorbing Fluorene Derivatives  
 AUTHOR(S): Belfield, Kevin D.; Morales, Alma R.; Kang, Bong-Soo; Hales, Joel M.; Hagan, David J.; Van Stryland, Eric W.; Chapela, Victor M.; Percino, Judith  
 CORPORATE SOURCE: Department of Chemistry and College of Optics and Photonics: CREOL FPCE, University of Central Florida, Orlando, FL, 32816, USA  
 SOURCE: Chemistry of Materials (2004), 16(23), 4634-4641  
 CODEN: CMATEX; ISSN: 0897-4756  
 PUBLISHER: American Chemical Society  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 OTHER SOURCE(S): CASREACT 141:313812

AB The synthesis of a series of four new compds. containing fluorenyl chromophores is presented, along with the results of spectroscopic and photochem. studies aimed at understanding the two-photon absorption properties and energetics of their electronically excited states. The mol. structures of the compds. were systematically varied to allow comparison of mols. possessing high and low mol. symmetry, short and long alkyl chains, and a fluorenyl conjugated  $\pi$ -system. Solvent-dependent absorption and emission were investigated along with  $\pi$ -conjugation length. Preliminary measurements of two-photon absorption (2PA) using a two-photon fluorescence method indicate that these chromophores exhibit high two-photon absorptivity. A sym. mol. (3), possessing a relatively large  $\pi$ -conjugated system, flanked on either side by electron-withdrawing groups (benzothiazole), exhibited a peak 2PA cross section ( $\delta$ ) of  $6000 + 10\text{-}50 \text{ cm}^4 \text{ s photon}^{-1} \text{ mol}^{-1}$  at 600 nm. Excitation anisotropy studies revealed the position of the  $S_0 \rightarrow S_1$  and  $S_0 \rightarrow S_2$  electronic transitions. Consistent with quantum mech. selection rules, the two-photon allowed transition ( $S_0 \rightarrow S_2$ ) was dominant.

IT 745079-44-1P  
 RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (synthesis, characterization, and optical spectra of two-photon-absorbing fluorene derivs.)  
 RN 745079-44-1 CAPLUS  
 CN 9H-Fluoren-2-amine, 7-[2-[4-(2-benzothiazolyl)phenyl]ethenyl]-9,9-didecyl-N,N-diphenyl- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 41 THERE ARE 41 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 27 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2004:360279 CAPLUS  
 DOCUMENT NUMBER: 140:392334  
 TITLE: Two-photon responsive chromophores containing electron accepting core units  
 INVENTOR(S): Kannan, Ramamurthi; Tan, Loon-seng; Reinhardt, Bruce A.; Vaia, Richard A.  
 PATENT ASSIGNEE(S): The United States of America as Represented by the Secretary of the Air Force, USA  
 SOURCE: U.S., 6 pp.  
 CODEN: USXXAM  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

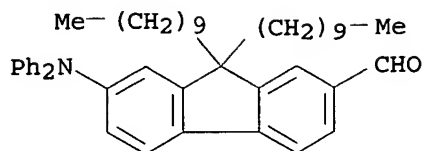
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6730793	B1	20040504	US 2002-171566	20020613
PRIORITY APPLN. INFO.:			US 2002-171566	20020613
OTHER SOURCE(S):	MARPAT 140:392334			
GI				

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB Chromophores are described by the general formula Q-(-L-Z)<sub>x</sub> (x = 2 or 3; Q is selected from I, II, III, IV, V, and VI; L = VII; R = C1-20 alkyl group; and Z = VIII or IX). The chromophores may exhibit high two-photon absorptions. Thus, 2,5-Bis(7-carbazol-9-yl-9,9-didecylfluoren-2-yl)-1,3-thiazolo(5,4d)1,3-thiazole was prepared and exhibited β = 2.8 cm/GW at 0.2 mol/L.

IT 348631-00-5P  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (chromophores with high two-photon absorptions)

RN 348631-00-5 CAPLUS  
 CN 9H-Fluorene-2-carboxaldehyde, 9,9-didecyl-7-(diphenylamino)- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 28 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:316280 CAPLUS

DOCUMENT NUMBER: 141:23125

TITLE: Two-photon absorption cross-sections of fluorene

derivatives with cationic substituents

AUTHOR(S): Kawamata, Jun; Akiba, Masaharu; Tani, Takeharu;

Harada, Akinori; Inagaki, Yoshio

CORPORATE SOURCE: Department of Chemistry and Earth Sciences, Faculty of Science, Yamaguchi University, Yamaguchi, 753-8512, Japan

SOURCE: Chemistry Letters (2004), 33(4), 448-449

CODEN: CMLTAG; ISSN: 0366-7022

PUBLISHER: Chemical Society of Japan

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Two novel fluorene derivs. having cationic substituents were synthesized. Two-photon absorption (TPA) properties of the derivs. were evaluated using the fluorescence-based technique with a femtosecond pulse emitted from a Ti:sapphire laser. Maximum TPA cross-sections of the compds. were estimated from the TPA spectra. TPA cross-sections of compds. with cationic substituents were found to be significantly larger than those of the compds. without cationic substituents.

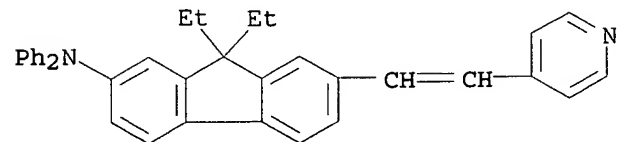
IT 197969-56-5 437713-15-0

RL: PRP (Properties)

(two-photon absorption cross-sections of fluorene derivs. with cationic substituents)

RN 197969-56-5 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



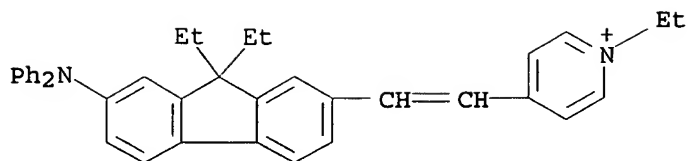
RN 437713-15-0 CAPLUS

CN Pyridinium, 4-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]-1-ethyl-, perchlorate (9CI) (CA INDEX NAME)

CM 1

CRN 437713-14-9

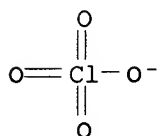
CMF C38 H37 N2



CM 2

CRN 14797-73-0

CMF Cl 04



REFERENCE COUNT: 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 29 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:210510 CAPLUS

DOCUMENT NUMBER: 140:414514

TITLE: Degenerate two-photon-absorption spectral studies of highly two-photon active organic chromophores  
 AUTHOR(S): He, Guang S.; Lin, Tzu-Chau; Dai, Jianming; Prasad, Paras N.; Kannan, Ramamurthi; Dombroskie, Ann G.; Vaia, Richard A.; Tan, Loon-Seng

CORPORATE SOURCE: Photonics and Biophotonics, Institute for Lasers, State University of New York at Buffalo, Buffalo, NY, 14260-3000, USA

SOURCE: Journal of Chemical Physics (2004), 120(11), 5275-5284  
 CODEN: JCPSA6; ISSN: 0021-9606

PUBLISHER: American Institute of Physics

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Degenerate two-photon absorption (TPA) spectral properties of five AFX chromophore solns. have been studied using a single and spectrally dispersed sub-picosecond white-light continuum beam. In a specially designed optical configuration, optical pathways inside the sample solution for different spectral components of the focused continuum beam were spatially separated from each other. Thus, the nondegenerate TPA processes coming from different spectral components can be eliminated, and the direct nonlinear absorption spectrum attributed to degenerate TPA processes can be readily obtained. Using this new technique, the complete TPA spectra for these five highly two-photon-active compds. (AF-380, AF-350, AF-295, AF-270, and AF-50) were obtained in the spectral range from 600 to 950 nm on an absolute scale of TPA cross section. The relationship between the mol. structures and their TPA spectral behaviors are discussed. In general the measured TPA spectra are not identical with the linear absorption spectra on the scale of absorbed photon(s) energy. Moreover, for some sample (such as AF-380), the TPA spectrum is totally different from the linear spectrum, which implies the difference of mol. transition pathways and selection rules for one- and two-photon excitation



processes. At high excitation intensity levels ( $\geq 15$  GW/cm<sup>2</sup>), the saturation behavior of TPA transition can be observed obviously in AF-350 and AF-380 solns. that exhibit much higher nonlinear absorptivity than the other chromophores investigated.

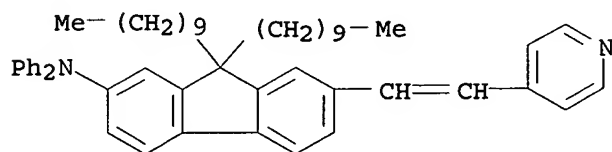
IT 191667-13-7, AF 50

RL: PRP (Properties)

(degenerate two-photon-absorption spectral studies of highly two-photon active dialkylfluorene-based chromophores)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



REFERENCE COUNT: 73 THERE ARE 73 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 30 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:165018 CAPLUS

DOCUMENT NUMBER: 141:233070

TITLE: Photostability of a series of two-photon absorbing fluorene derivatives

AUTHOR(S): Belfield, Kevin D.; Bondar, Mykhailo V.; Przhonska, Olga V.; Schafer, Katherine J.

CORPORATE SOURCE: Department of Chemistry and CREOL/School of Optics, University of Central Florida, Orlando, FL, 32816-2366, USA

SOURCE: Journal of Photochemistry and Photobiology, A: Chemistry (2004), 162(2-3), 489-496  
CODEN: JPPCEJ; ISSN: 1010-6030

PUBLISHER: Elsevier Science B.V.

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The photochem. stability of a series of two-photon absorbing (TPA) fluorene derivs. was investigated in air- and N<sub>2</sub>-saturated acetonitrile (ACN) at room temperature. The quantum yields of the photoreactions,  $\Phi$ , were determined at various concns. of the fluorene derivs., oxygen concentration of the solvent,

and irradiation wavelength. The absorption and fluorescence spectra of the photoproducts, corresponding to different excitation conditions, were analyzed. Photooxidn. and electron transfer processes are proposed as photobleaching mechanisms for the fluorene derivs. in ACN. The relatively low photochem. quantum yields ( $\Phi$  .apprx.10<sup>-4</sup>) make the derivs. particularly promising for linear and nonlinear optical applications.

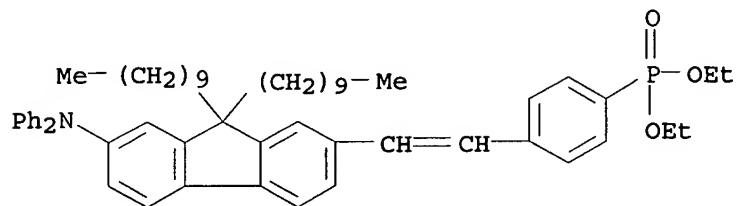
IT 507271-13-8

RL: CPS (Chemical process); PEP (Physical, engineering or chemical process); PRP (Properties); RCT (Reactant); PROC (Process); RACT (Reactant or reagent)

(photochem. properties of two-photon absorbing fluorene derivs. in acetonitrile solution as function of concentration and oxygen content and irradiation wavelength)

RN 507271-13-8 CAPLUS

CN Phosphonic acid, [4-[2-[9,9-didecyl-7-(diphenylamino)-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)



REFERENCE COUNT: 19 THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 31 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2004:76492 CAPLUS  
 DOCUMENT NUMBER: 140:136180  
 TITLE: Amorphous polyphenol derivatives with good heat resistance and organic electroluminescent devices  
 INVENTOR(S): Fukuoka, Naohiko; Tagami, Sanae; Fujiwara, Toru; Shionoya, Hidehiko  
 PATENT ASSIGNEE(S): Chemipro Kasei Ltd., Japan  
 SOURCE: Jpn. Kokai Tokkyo Koho, 67 pp.  
 CODEN: JKXXAF  
 DOCUMENT TYPE: Patent  
 LANGUAGE: Japanese  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2004026757	A2	20040129	JP 2002-188237	20020627
PRIORITY APPLN. INFO.:			JP 2002-188237	20020627
OTHER SOURCE(S):	MARPAT 140:136180			
GI				

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB The derivs. are I [Q = A, B, etc.; R1, R20, R21 = H, (halo)alkyl, cycloalkyl, etc.; R2-R19 = H, halo, (halo)alkyl, etc.; M = C, D; Ar1 = arylene, oxydiaryldiyl; R22 = H, (cyclo)alkyl, aryl; R23-R26 = H, alkyl, alkoxy, aryl; R27-R30 = H, alkyl, alkoxy, aralkyl, etc.; h, m, n = 1-3; j, k, p = 1-4]. Emitter or hole-transport layers of the devices are easily manufactured by solution casting of the derivs. without polymeric binders.

IT 648908-22-9P 648908-23-0P

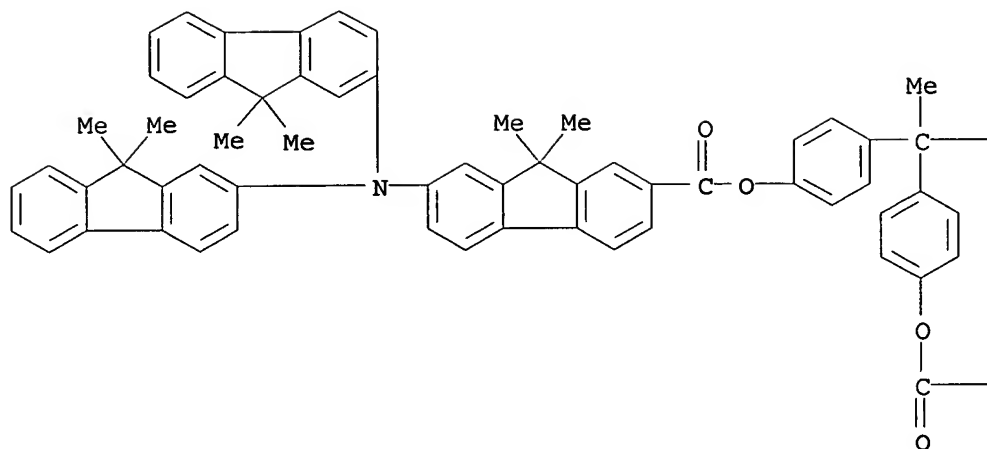
RL: DEV (Device component use); IMF (Industrial manufacture); PREP (Preparation); USES (Uses)

(heat-resistant amorphous polyphenol derivs. suitable for solution casting for manufacture of organic electroluminescent devices)

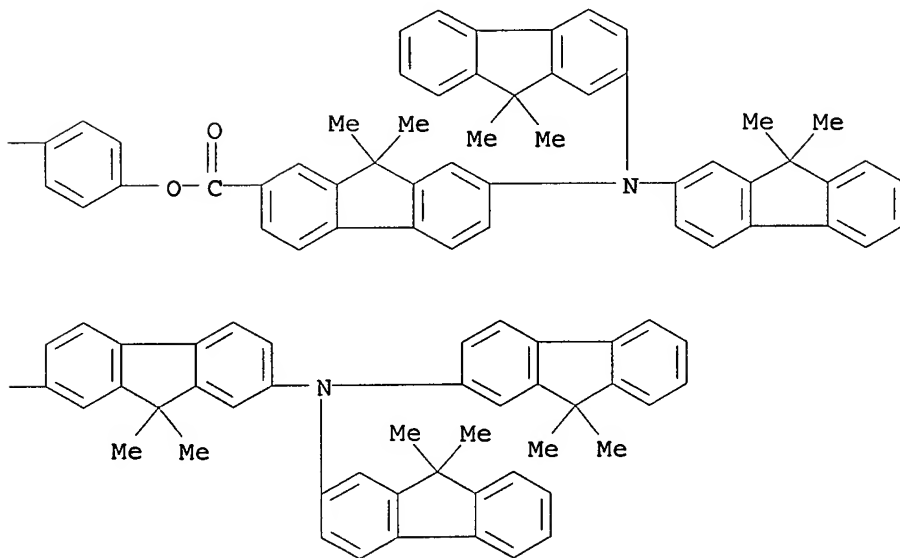
RN 648908-22-9 CAPLUS

CN 9H-Fluorene-2-carboxylic acid, 7-[bis(9,9-dimethyl-9H-fluoren-2-yl)amino]-9,9-dimethyl-, ethylidynetri-4,1-phenylene ester (9CI) (CA INDEX NAME)

PAGE 1-A

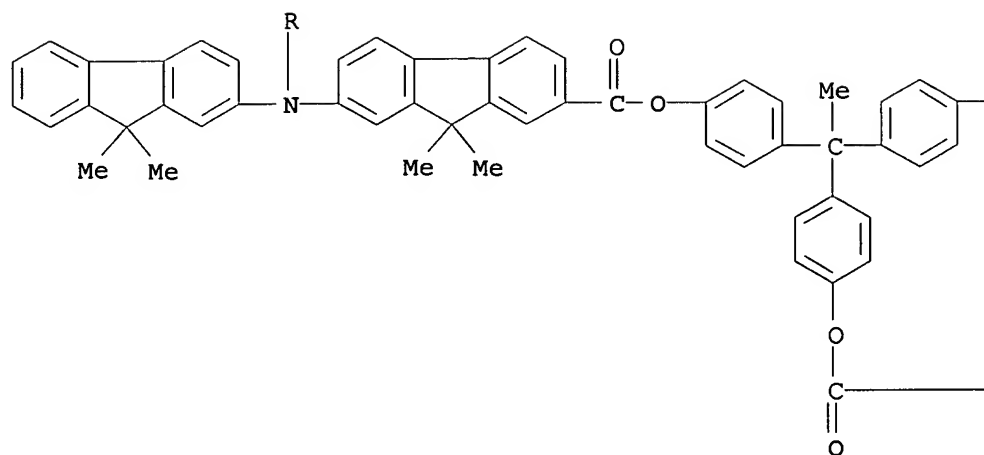


PAGE 1-B

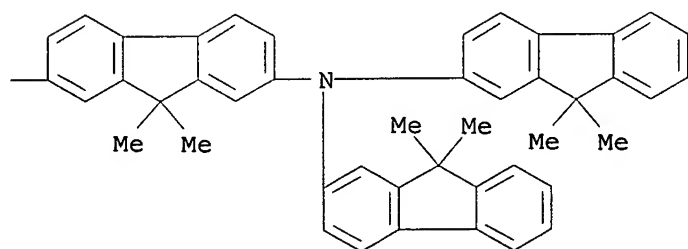
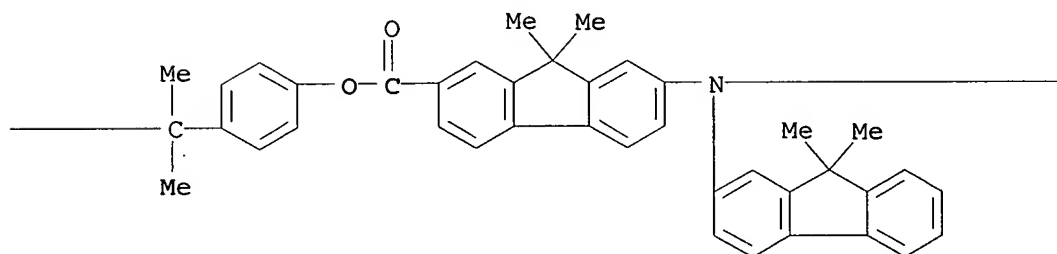


RN 648908-23-0 CAPLUS  
 CN 9H-Fluorene-2-carboxylic acid, 7-[bis(9,9-dimethyl-9H-fluorene-2-yl)amino] -  
 9,9-dimethyl-, [1-[4-[1-[4-[[[7-[bis(9,9-dimethyl-9H-fluorene-2-yl)amino] -  
 9,9-dimethyl-9H-fluorene-2-yl]carbonyl]oxy]phenyl]-1-  
 methylethyl]phenyl]ethylidene]di-4,1-phenylene ester (9CI) (CA INDEX  
 NAME)

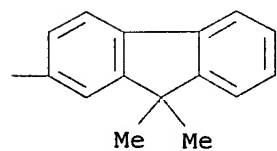
PAGE 1-A

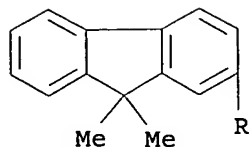


PAGE 1-B



PAGE 1-C





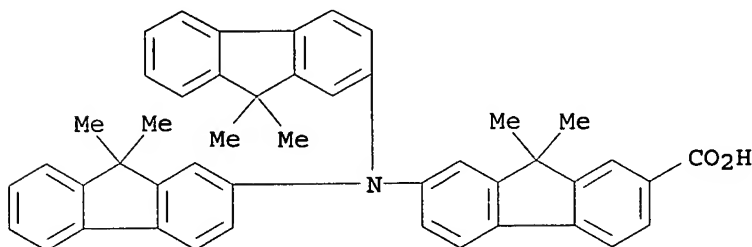
IT 648908-21-8

RL: RCT (Reactant); RACT (Reactant or reagent)

(heat-resistant amorphous polyphenol derivs. suitable for solution casting for manufacture of organic electroluminescent devices)

RN 648908-21-8 CAPLUS

CN 9H-Fluorene-2-carboxylic acid, 7-[bis(9,9-dimethyl-9H-fluorene-2-yl)amino]-9,9-dimethyl- (9CI) (CA INDEX NAME)



L6 ANSWER 32 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:60449 CAPLUS

DOCUMENT NUMBER: 140:128164

TITLE: Preparation and laser irradiation of fullerene derivatives for generation of singlet oxygen useful for treatment of tumors

INVENTOR(S): Chiang, Long Y.; Tan, Loon-Seng

PATENT ASSIGNEE(S): USA

SOURCE: PCT Int. Appl., 28 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004007426	A1	20040122	WO 2003-US22265	20030716
W:	AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2003249294	A1	20040202	AU 2003-249294	20030716

US 2005191229	A1	20050901	US 2003-620839	20030716
PRIORITY APPLN. INFO.:			US 2002-396782P	P 20020717
			WO 2003-US22265	W 20030716
OTHER SOURCE(S):	MARPAT 140:128164			
GI				

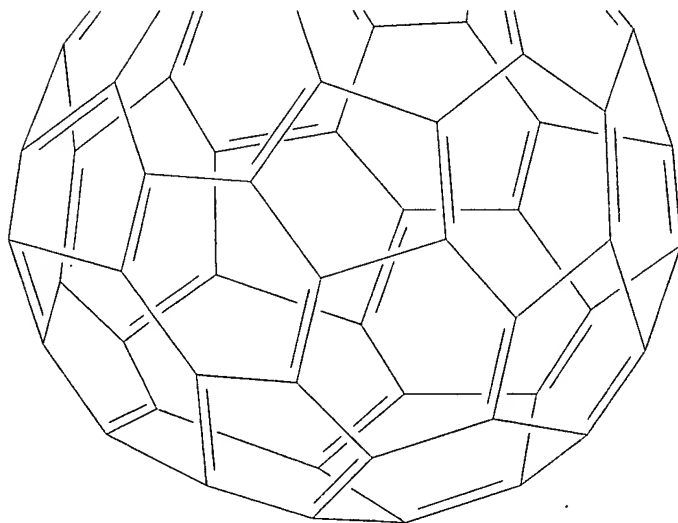
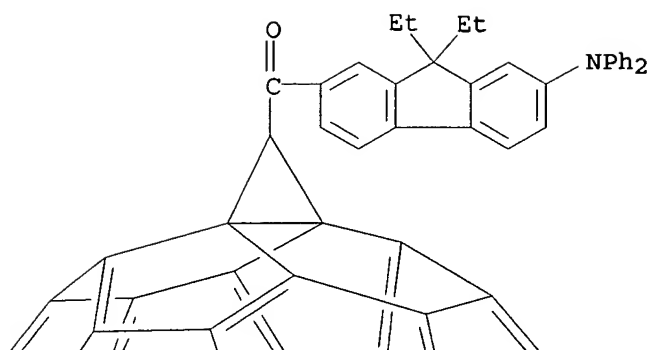
\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB The invention relates to fullerene compds. of formula I [wherein: R1, R2 = 3,4-disubstituted phenyl; R3, R4, R5 = (un)substituted aryl(oxy/thio), alkyl, alkylaryl, etc.; V = (hetero)aryl; W = O, C(CN)2, etc.; F = fullerene core; R = OH, NH2; E is represented by wide variety of substituents; p, q = 0-20; n = 1-10], laser irradiation of which generates singlet oxygen, useful for inhibition of tumor growth. Also disclosed are pharmaceutical compns. containing one of the compds. described above. For instance, compound II [X-Y is a 6/6 C-C bond of fullerene C60] was prepared via methoxyethylation of 2-bromofluorene, amination of the obtained 2-bromo-9,9-bis(2-methoxyethyl)fluorene by diphenylamine, and subsequent bromoacetylation by BrCH2C(O)Br, addition of fullerene C60 to the obtained fluorene derivative III, bis-O-demethylation, and etherification by 1,4-butane sultone. In a laser irradiation experiment using another compound I, an unexpectedly large amount of singlet oxygen, and oxygenated radicals generated from fullerene derivs. or composites, were detected. Both in vitro and in vivo studies of I showed inhibition of tumor cell growth (no data).

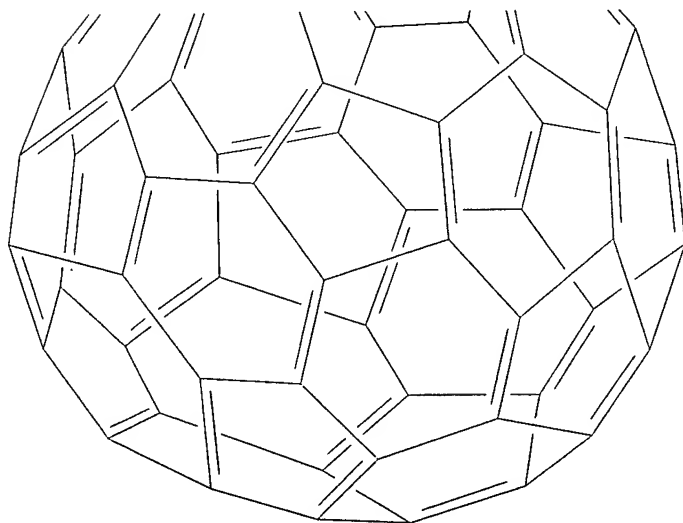
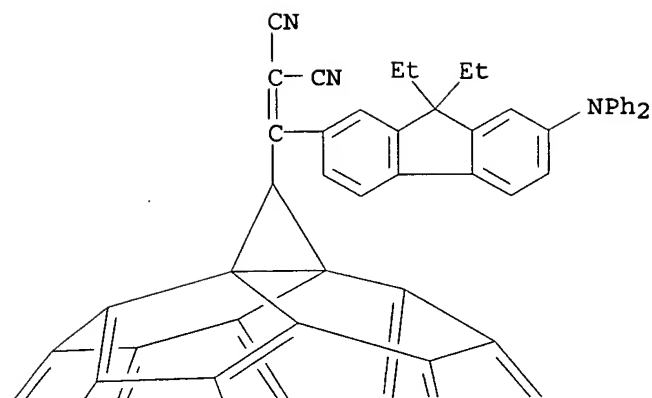
IT 486998-58-7P, 7-[1,2-Dihydro-1,2-methano(60)fullerene-61-carbonyl]-9,9-diethyl-2-diphenylaminofluorene 649724-43-6P  
 RL: PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (intermediate; preparation and laser irradiation of fullerene derivs. for generation of singlet oxygen useful for inhibition of tumor growth)

RN 486998-58-7 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



RN 649724-43-6 CAPLUS  
 CN Propanedinitrile, [3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]methylene]- (9CI) (CA INDEX NAME)



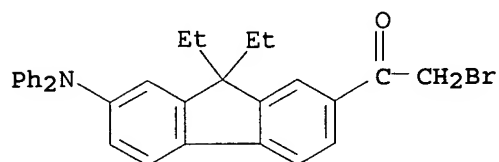
IT 486998-57-6P, 7-Bromoacetyl-9,9-diethyl-2-diphenylaminofluorene  
 649724-47-0P, 7-Bromoacetyl-9,9-bis(2-methoxyethyl)-2-  
 diphenylaminofluorene 649724-49-2P, 7-[1,2-Dihydro-1,2-  
 methano(60) fullerene-61-carbonyl]-9,9-bis(2-methoxyethyl)-2-  
 diphenylaminofluorene 649724-50-5P, 7-[1,2-Dihydro-1,2-  
 methano(60) fullerene-61-carbonyl]-9,9-bis(2-hydroxyethyl)-2-  
 diphenylaminofluorene



RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(intermediate; preparation and laser irradiation of fullerene derivs. for generation of singlet oxygen useful for inhibition of tumor growth)

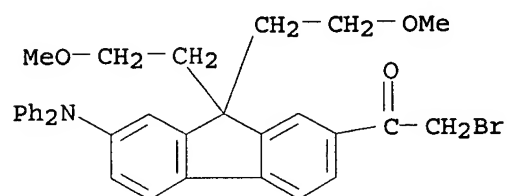
RN 486998-57-6 CAPLUS

CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI)  
(CA INDEX NAME)



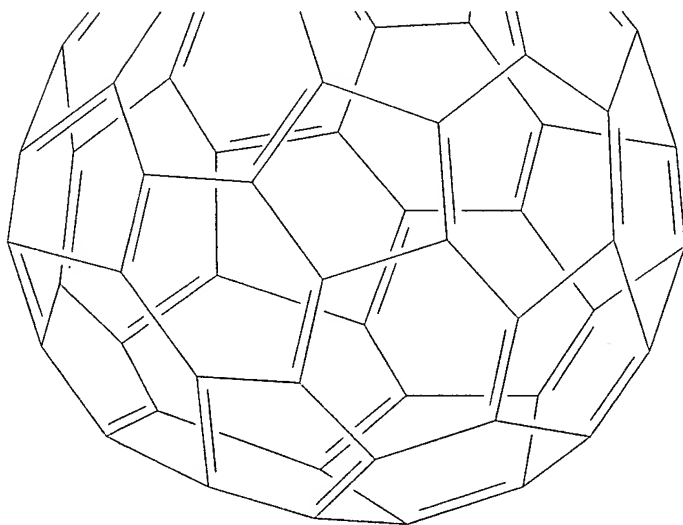
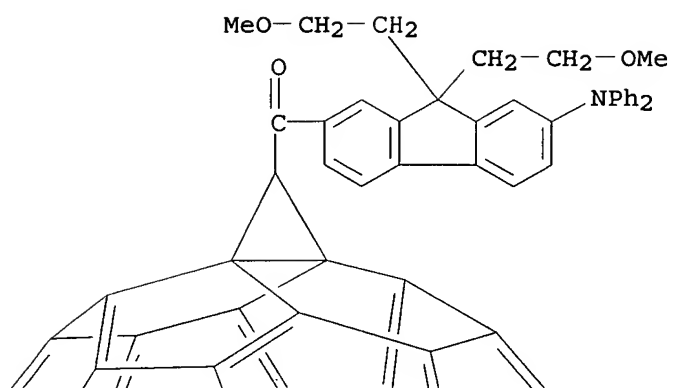
RN 649724-47-0 CAPLUS

CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



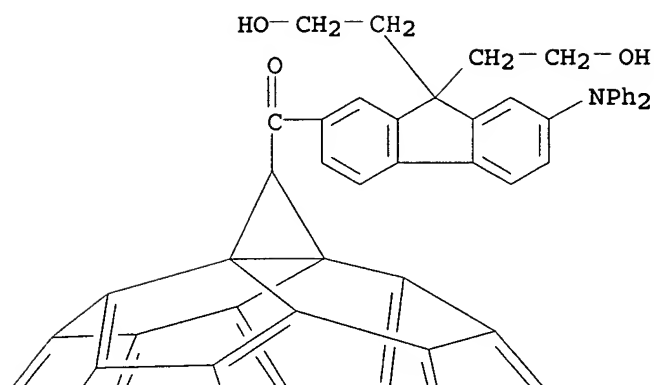
RN 649724-49-2 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-bis(2-methoxyethyl)-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)

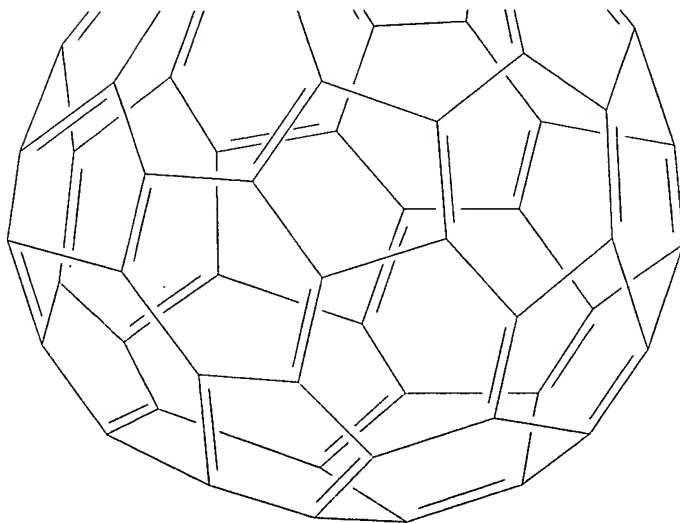


RN 649724-50-5 CAPLUS  
 CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-bis(2-hydroxyethyl)-9H-fluoren-2-yl]-(9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 2-A

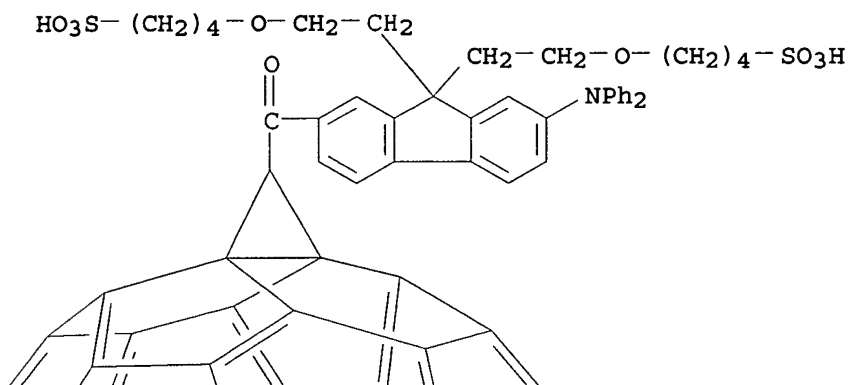


IT 649724-51-6P, 7-[1,2-Dihydro-1,2-methano(60) fullerene-61-carbonyl] -  
 9,9-bis[2-(4-sulfobutoxy)ethyl]-2-diphenylaminofluorene  
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU  
 (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES  
 (Uses)  
 (preparation and laser irradiation of fullerene derivs. for generation of  
 singlet oxygen useful for inhibition of tumor growth)

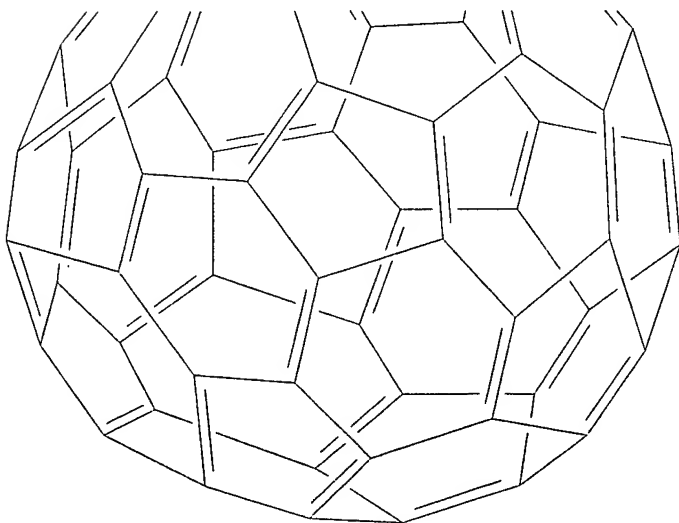
RN 649724-51-6 CAPLUS

CN 1-Butanesulfonic acid, [[2-(3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-ylcarbonyl)-7-(diphenylamino)-9H-fluoren-9-ylidene]bis(2,1-ethanediylloxy)]bis- (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 2-A



REFERENCE COUNT:

2

THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

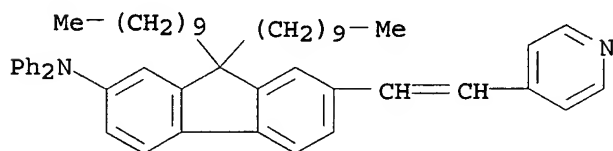
L6 ANSWER 33 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:10215 CAPLUS  
 DOCUMENT NUMBER: 141:30387  
 TITLE: Two-photon excitation of dyes in a polymer matrix by femtosecond pulses from a Ti:sapphire laser  
 AUTHOR(S): Meshalkin, Yu. P.; Svetlichnyi, V. A.; Reznichenko, A. V.; Myachin, A. Yu.; Bakhareva, S. S.; Dolotov, S. M.; Kopylova, T. N.; Ponomarenko, E. P.  
 CORPORATE SOURCE: Center for Photodynamic Research, Research Institute of Physiology, Siberian Branch, Russian Academy of Medical Sciences, Novosibirsk, 630117, Russia  
 SOURCE: Quantum Electronics (2003), 33(9), 803-806  
 CODEN: QUELEZ; ISSN: 1063-7818  
 PUBLISHER: Turpion Ltd.  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English

AB Two-photon fluorescence was observed for 18 organic dyes in a PMMA matrix excited by a fs Ti:Al<sub>2</sub>O<sub>3</sub> laser. The product of the cross section for 2-photon absorption by the quantum yield of fluorescence (2-photon fluorescence cross section) is estimated by comparing it with fluorescence of Rhodamine 6G in EtOH. Using this parameter, dyes are selected that exhibit the most intense fluorescence in PMMA and their concns. in PMMA are optimized. Coumarin and rhodamine dyes in polymer matrixes are proposed for using as visualizers of fs radiation of a Ti:Al<sub>2</sub>O<sub>3</sub> laser and as detectors in self-triggering systems.

IT 191667-13-7  
 RL: MOA (Modifier or additive use); PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); PROC (Process); USES (Uses)  
 (two-photon excitation by fs pulses from titanium-doped alumina laser of PMMA matrix containing)

RN 191667-13-7 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



REFERENCE COUNT: 20 THERE ARE 20 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 34 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:891017 CAPLUS  
 DOCUMENT NUMBER: 141:196943  
 TITLE: Nonlinear optical spectroscopic characterization of a series of fluorene derivatives  
 AUTHOR(S): Hales, Joel M.; Schafer, Katherine J.; Morales, Alma M.; Belfield, Kevin D.; Hagan, David J.; Van Stryland, Eric W.  
 CORPORATE SOURCE: School of Optics/CREOL, Univ. of Central Florida, Orlando, FL, 32816, USA  
 SOURCE: Proceedings of SPIE-The International Society for Optical Engineering (2003), 5211(Nonlinear Optical Transmission and Multiphoton Processes in Organics),

21-30

CODEN: PSISDG; ISSN: 0277-786X

PUBLISHER:

SPIE-The International Society for Optical Engineering

DOCUMENT TYPE:

Journal

LANGUAGE:

English

AB The authors have performed nonlinear spectroscopic measurements to study the chemical structure/nonlinear optical property relations for a set of alkyl fluorene derivs. The characterization method the authors used is a femtosecond white-light continuum (WLC) pump-probe spectrometer that can rapidly characterize an organic samples nondegenerate two-photon absorption (2PA) spectrum. The nature of these expts. requires sophisticated data anal. In particular, the relative group velocity mismatch between the pump and probe, which are at different frequencies, makes these pulses walk through each other within the thickness of the sample. For widely different frequencies, this can severely diminish the 2PA signal strength. However, given careful anal., the authors found good agreement with known semiconductor samples. Confidence in this method has allowed the authors to study the effects of solvent effect, electron-withdrawing character, conjugation length, and symmetry on the two-photon absorbing properties of these mols. The authors found an optimum solvent polarity as well as electron-withdrawing character which serves to maximize the strength of the 2PA in these materials. Different synthesis avenues provided the authors with two different methods of extending the conjugation length that increases the nonlinearity as well. Finally, studies of mols. with disparate symmetry have allowed the authors to identify the symmetry of the excited states. The authors present the 1st exptl. study of the intermediate state resonance enhancement of nondegenerate 2PA in organic mols. Using a simplified sum-over-states expression, the authors make comparisons between experiment and theory.

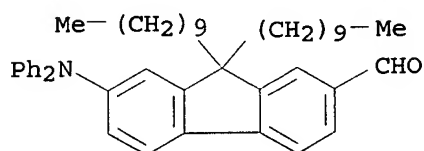
IT 348631-00-5 738584-52-6 738584-53-7

RL: PRP (Properties)

(nonlinear optical spectroscopic characterization of a series of fluorene derivs.)

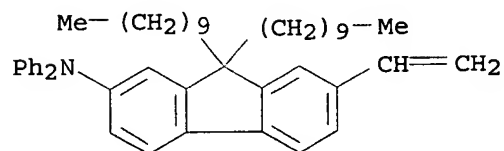
RN 348631-00-5 CAPLUS

CN 9H-Fluorene-2-carboxaldehyde, 9,9-didecyl-7-(diphenylamino)- (9CI) (CA INDEX NAME)



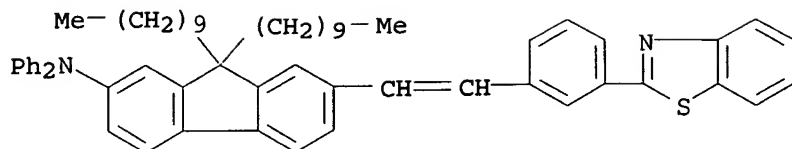
RN 738584-52-6 CAPLUS

CN 9H-Fluorene-2-amine, 9,9-didecyl-7-ethenyl-N,N-diphenyl- (9CI) (CA INDEX NAME)



RN 738584-53-7 CAPLUS

CN 9H-Fluorene-2-amine, 7-[2-[3-(2-benzothiazolyl)phenyl]ethenyl]-9,9-didecyl-N,N-diphenyl- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 35 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:702370 CAPLUS

DOCUMENT NUMBER: 140:41640

TITLE: Influence of electron-acceptor strength on the resonant two-photon absorption cross sections of diphenylaminofluorene-based chromophores

AUTHOR(S): Guo, Jing-Dong; Wang, Chuan-Kui; Luo, Yi; Agren, Hans  
CORPORATE SOURCE: Theoretical Chemistry, SCFAB, Royal Institute of Technology, Stockholm, S-10691, Swed.

SOURCE: Physical Chemistry Chemical Physics (2003), 5(18), 3869-3873

CODEN: PPCPFQ; ISSN: 1463-9076

PUBLISHER: Royal Society of Chemistry

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Resonant two-photon absorption (TPA) cross sections of a series of diphenylaminofluorene-based chromophores with various electron acceptors are predicted using the RPA and using hybrid d. functional theory implemented for a two-state model. A comparison of the two methods indicates that the two-state model is adequate for describing the TPA cross sections of all asym. charge-transfer systems under investigation. It is demonstrated that the inclusion of electron correlation can drastically increase the absolute values of the TPA cross sections, but that it has negligible effects on the relative order of the TPA activity of the mols.

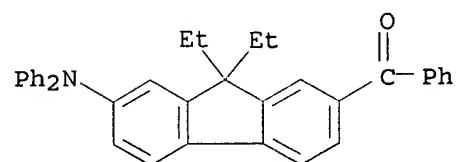
IT 349123-45-1, AF-370

RL: PRP (Properties)

(influence of electron-acceptor strength on resonant two-photon absorption cross sections of diphenylaminofluorene-based chromophores)

RN 349123-45-1 CAPLUS

CN Methanone, [7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]phenyl- (9CI)  
(CA INDEX NAME)



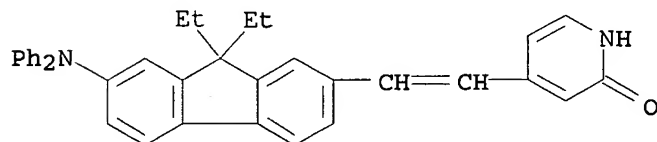
REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 36 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

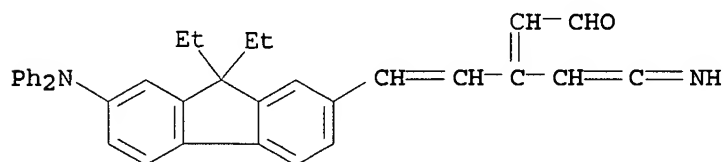
ACCESSION NUMBER: 2003:686696 CAPLUS

DOCUMENT NUMBER: 140:50152

TITLE: A theoretical study of the photooxidation of a fluorene-based two-photon chromophore  
 AUTHOR(S): Guo, Jing-Dong; Luo, Yi  
 CORPORATE SOURCE: Laboratory of Theoretical Chemistry, Royal Institute of Technology, Stockholm, S-10691, Swed.  
 SOURCE: THEOCHEM (2003), 635, 1-9  
 CODEN: THEODJ; ISSN: 0166-1280  
 PUBLISHER: Elsevier Science B.V.  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 AB The IR and optical absorption spectra of a fluorene-based two-photon chromophore AF50 (N,N-diphenyl-7-2-(4-pyridinyl)-ethenyl-9,9-di-n-decyl-9H-fluorene-2-amine) and its photooxidn. products after UV exposure have been calculated using hybrid d. functional theory. The authors calcns. have provided very useful information about the nature of the read/write process of AF50 and its possible final products.  
 IT 634908-90-0 634908-91-1 634908-92-2  
 634908-93-3 634908-94-4 634908-95-5  
 634908-96-6 634908-97-7 634908-98-8  
 RL: CPS (Chemical process); FMU (Formation, unclassified); PEP (Physical, engineering or chemical process); PRP (Properties); FORM (Formation, nonpreparative); PROC (Process)  
 (photoproduct; theor. study of photooxidn. of fluorene-based two-photon chromophore used in optical memory devices)  
 RN 634908-90-0 CAPLUS  
 CN 2(1H)-Pyridinone, 4-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]- (9CI) (CA INDEX NAME)

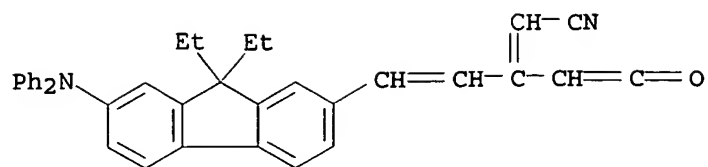


RN 634908-91-1 CAPLUS  
 CN 2,4-Pentadienal, 3-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]-5-imino- (9CI) (CA INDEX NAME)

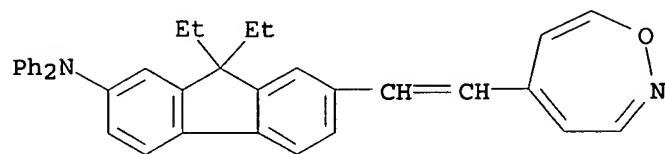


RN 634908-92-2 CAPLUS  
 CN 2,4-Pentadienenitrile, 3-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]-5-oxo- (9CI) (CA INDEX NAME)

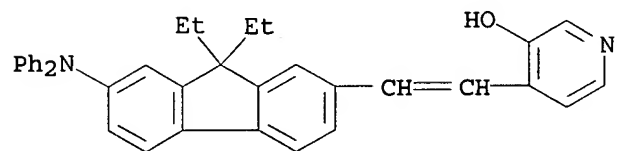




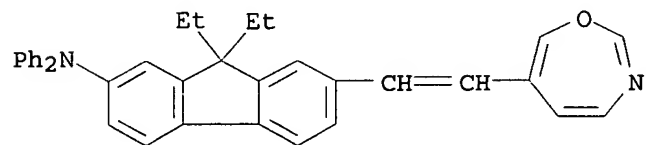
RN 634908-93-3 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-7-[2-(1,2-oxazepin-5-yl)ethenyl]-N,N-diphenyl- (9CI) (CA INDEX NAME)



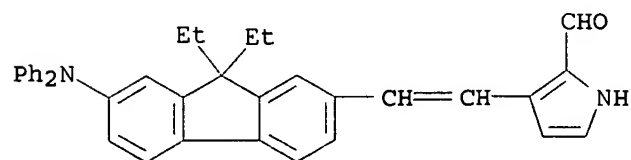
RN 634908-94-4 CAPLUS  
 CN 3-Pyridinol, 4-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]- (9CI) (CA INDEX NAME)



RN 634908-95-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-7-[2-(1,3-oxazepin-6-yl)ethenyl]-N,N-diphenyl- (9CI) (CA INDEX NAME)

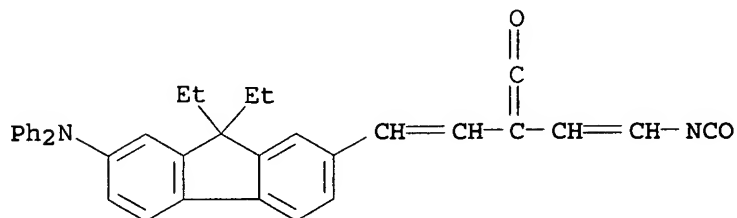


RN 634908-96-6 CAPLUS  
 CN 1H-Pyrrole-2-carboxaldehyde, 3-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]- (9CI) (CA INDEX NAME)



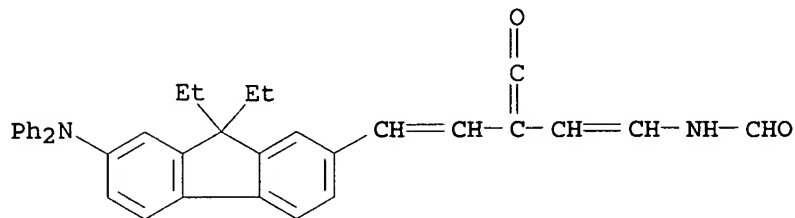
RN 634908-97-7 CAPLUS

CN 1,3-Butadien-1-one, 2-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]-4-isocyanato- (9CI) (CA INDEX NAME)



RN 634908-98-8 CAPLUS

CN Formamide, N-[3-carbonyl-5-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]-1,4-pentadienyl]- (9CI) (CA INDEX NAME)



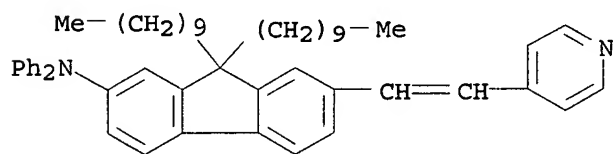
IT 191667-13-7, AF50 197969-56-5

RL: CPS (Chemical process); PEP (Physical, engineering or chemical process); PRP (Properties); RCT (Reactant); PROC (Process); RACT (Reactant or reagent)

(theor. study of photooxidn. of fluorene-based two-photon chromophore used in optical memory devices)

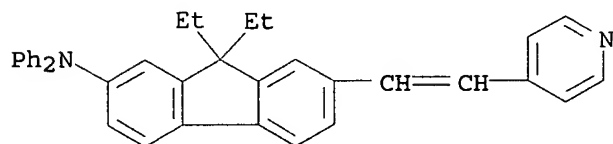
RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



RN 197969-56-5 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 37 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:600631 CAPLUS

DOCUMENT NUMBER: 139:314360

TITLE: Inter- and Intramolecular Photoinduced  
Electron-Transfer Processes between C60 and  
Diphenylaminofluorene in SolutionsAUTHOR(S): Luo, Hongxia; Fujitsuka, Mamoru; Araki, Yasuyuki; Ito,  
Osamu; Padmawar, Prashant; Chiang, Long Y.CORPORATE SOURCE: Institute of Multidisciplinary Research for Advanced  
Materials, Tohoku University, Aoba, Sendai, 980-8577,  
JapanSOURCE: Journal of Physical Chemistry B (2003), 107(35),  
9312-9318

CODEN: JPCBPK; ISSN: 1520-6106

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

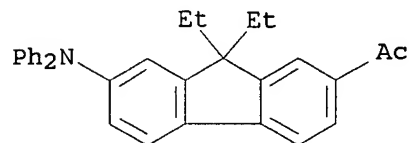
AB Intramol. photoinduced charge-separation and charge-recombination processes of a covalently bonded buckminsterfullerene-diphenylaminofluorene (C60-DPAF) dyad, in which the diphenylaminofluorene moiety is connected to C60 via a ketone group, have been investigated. Quenching of the fluorescence intensities and the observed short fluorescence lifetimes of the C60 moiety of the dyad in benzonitrile (PhCN) and DMF indicated that charge separation takes place via the singlet excited state of the C60 moiety at a fast rate with high efficiency. On the basis of the nanosecond transient absorption spectra, formation of the radical ion pair  $C60^{\bullet-}$ -DPAF $^{\bullet+}$  was confirmed in DMF; the radical ion pair decays with a lifetime of 150 ns. From the temperature dependence of the charge-recombination rate constants, the reorganization energy was evaluated to be 0.81 eV in DMF, which is reasonably small, characteristic of fullerene derivs. On the other hand, a mixture of DPAF and C60 showed intermol. electron transfer via the triplet state of C60 in polar solvents such as PhCN.

IT 612057-44-0

RL: CPS (Chemical process); FMU (Formation, unclassified); PEP (Physical, engineering or chemical process); PRP (Properties); FORM (Formation, nonpreparative); PROC (Process)  
(inter- and intramol. photoinduced electron-transfer processes between C60 and diphenylaminofluorene in solns.)

RN 612057-44-0 CAPLUS

CN Ethanone, 1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]-, radical ion(1+) (9CI) (CA INDEX NAME)

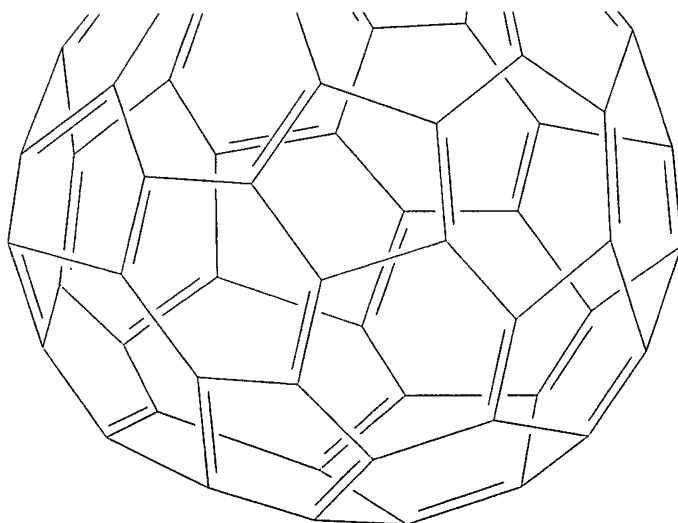
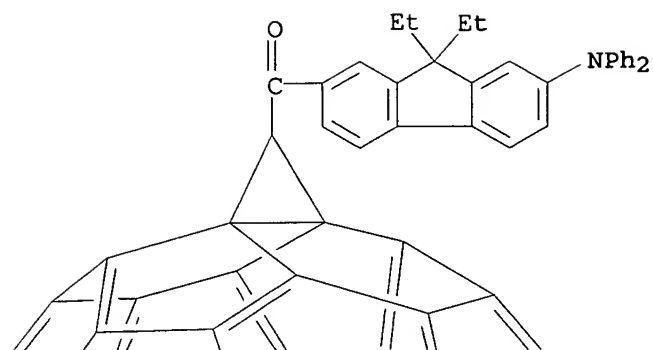


IT 486998-58-7 612057-42-8

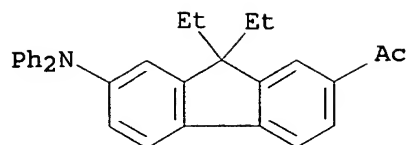
RL: CPS (Chemical process); PEP (Physical, engineering or chemical process); PRP (Properties); PROC (Process)  
(inter- and intramol. photoinduced electron-transfer processes between C60 and diphenylaminofluorene in solns.)

RN 486998-58-7 CAPLUS

CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



RN 612057-42-8 CAPLUS  
 CN Ethanone, 1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl] - (9CI) (CA  
 INDEX NAME)



REFERENCE COUNT: 34 THERE ARE 34 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 38 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:293006 CAPLUS

DOCUMENT NUMBER: 139:36200

TITLE: Chemical structure/nonlinear optical property

relations for fluorenyl ring system derivatives

AUTHOR(S): Hales, J.; Schafer, K. J.; Morales, A. M.; Belfield, K. D.; Hagan, D. J.; Van Stryland, E. W.

CORPORATE SOURCE: School of Optics/CREOL, University of Central Florida, Orlando, FL, 32816-2700, USA

SOURCE: Trends in Optics and Photonics (2002), 79(Nonlinear Optics), 369-371

CODEN: TOPRBS

PUBLISHER: Optical Society of America

DOCUMENT TYPE: Journal

LANGUAGE: English

AB We present initial work involving chemical-structure/nonlinear-optical (NLO) property relations for a set of fluorene derivs. This is achievable using our femtosecond white-light continuum pump-probe nonlinear spectrometer which can rapidly characterize a sample's two-photon absorption spectrum.

IT 540536-57-0

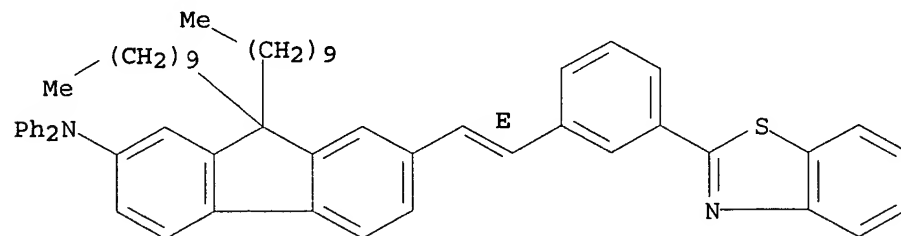
RL: PRP (Properties)

(chemical structure-nonlinear optical property relations for fluorenyl ring system derivs.)

RN 540536-57-0 CAPLUS

CN 9H-Fluoren-2-amine, 7-[(1E)-2-[3-(2-benzothiazolyl)phenyl]ethenyl]-9,9-didecyl-N,N-diphenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 39 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:71489 CAPLUS

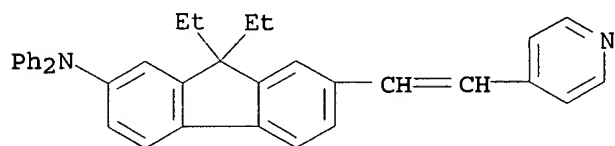
DOCUMENT NUMBER: 138:115140

TITLE: Two-photon absorption-type optical information recording media

INVENTOR(S): Inagaki, Yoshio; Akiba, Masaharu

PATENT ASSIGNEE(S): Fuji Photo Film Co., Ltd., Japan  
 SOURCE: Jpn. Kokai Tokkyo Koho, 13 pp.  
 CODEN: JKXXAF  
 DOCUMENT TYPE: Patent  
 LANGUAGE: Japanese  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

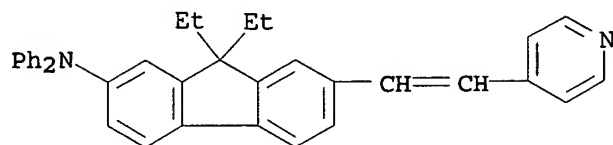
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2003029376	A2	20030129	JP 2001-216758	20010717
PRIORITY APPLN. INFO.:			JP 2001-216758	20010717
AB A highly sensitive recording medium, which requires relatively low power laser to provoke two-photon absorption, is characterized by containing a fluorescence quencher and a two-photon absorbing compound having a two-photon absorption cross-section >102 GM (1 GM = 1 + 10 <sup>-50</sup> cm <sup>4</sup> s mol. <sup>-1</sup> photon <sup>-1</sup> ).				
IT 197969-56-5				
RL: DEV (Device component use); USES (Uses) (two-photon absorbing compound in optical information recording media)				
RN 197969-56-5 CAPLUS				
CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)				



L6 ANSWER 40 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2003:56348 CAPLUS  
 DOCUMENT NUMBER: 138:114755  
 TITLE: Composite for two photon absorption  
 INVENTOR(S): Inagaki, Yoshio; Akiba, Masaharu; Harada, Akinori;  
 Tani, Takeharu  
 PATENT ASSIGNEE(S): Fuji Photo Film Co., Ltd., Japan  
 SOURCE: Jpn. Kokai Tokkyo Koho, 14 pp.  
 CODEN: JKXXAF  
 DOCUMENT TYPE: Patent  
 LANGUAGE: Japanese  
 FAMILY ACC. NUM. COUNT: 2  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2003020469	A2	20030124	JP 2001-206780	20010706
US 2003052311	A1	20030320	US 2002-188959	20020705
PRIORITY APPLN. INFO.:			JP 2001-206780	A 20010706
			JP 2001-212310	A 20010712
AB The invention refers to a composite with a two-photon absorption cross section > 102 Goppert-Mayers (GM) where 1 GM = 1 x 10 <sup>-50</sup> cm <sup>4</sup> s mol. <sup>-1</sup> photon <sup>-3</sup> in order to achieve two photon absorption using relatively low-powered lasers.				
IT 197969-56-5				
RL: DEV (Device component use); USES (Uses) (conjugated aryl amine and luminescent compound as composite for two				

photon absorption)  
 RN 197969-56-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



L6 ANSWER 41 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2002:924150 CAPLUS

DOCUMENT NUMBER: 138:287182

TITLE: Steady-State Spectroscopic and Fluorescence Lifetime Measurements of New Two-Photon Absorbing Fluorene Derivatives

AUTHOR(S): Belfield, Kevin D.; Bondar, Mikhailo V.; Przhonska, Olga V.; Schafer, Katherine J.

CORPORATE SOURCE: Department of Chemistry, University of Central Florida, Orlando, FL, USA

SOURCE: Journal of Fluorescence (2002), 12(3/4), 449-454

CODEN: JOFLEN; ISSN: 1053-0509

PUBLISHER: Kluwer Academic/Plenum Publishers

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Steady-state excitation anisotropy, lifetimes, and time-resolved emission spectra of new 2-photon absorbing fluorene derivs. [(7-benzothiazol-2-yl-9,9-didecylfluoren-2-yl)diphenylamine, 9,9-didecyl-2,7-bis(N,N-diphenylamino)fluorene, and {4-[2-(7-diphenylamino-9,9-diethylfluoren-2-yl)-vinyl]phenyl}phosphoric acid di-Et ester] were measured in aprotic solvents at room temperature. Excitation anisotropy spectra in viscous silicon oil allowed the determination of the spectral position of three electronic transitions S0 S1, S0 S2, S0 S3 (Si, i = 1, 2, 3 are the singlet electronic states) and the angles (.simeq. 30°) between absorption S0 S1 and emission S1 S0 dipole moments for the first electronic transition. Solvate relaxation processes in the first excited state of the investigated fluorene mols. affect the lifetimes of these states,  $\tau_1$ , so that exptl. values of  $\tau_1$  do not correspond to those calculated by Strickler and Berg theory. The influence of the mol. concentration on the fluorescence quantum yields and  $\tau_1$  have been investigated.

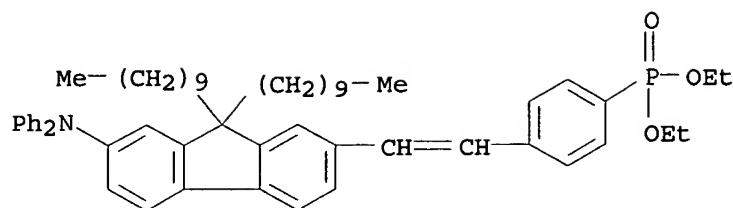
IT 507271-13-8

RL: PRP (Properties)

(steady-state spectroscopic and fluorescence lifetime measurements of new two-photon absorbing fluorene derivs.)

RN 507271-13-8 CAPLUS

CN Phosphonic acid, [4-[2-[9,9-didecyl-7-(diphenylamino)-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)



REFERENCE COUNT: 23 THERE ARE 23 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 42 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2002:647374 CAPLUS

DOCUMENT NUMBER: 138:106347

TITLE: Synthesis of C60-diphenylaminofluorene dyad with large 2PA cross-sections and efficient intramolecular two-photon energy transfer

AUTHOR(S): Chiang, Long Y.; Padmawar, Prashant A.; Canteenwala, Taizoon; Tan, Loon-Seng; He, Guang S.; Kannan, Ramamurthi; Vaia, Richard; Lin, Tzu-Chau; Zheng, Qingdong; Prasad, Paras N.

CORPORATE SOURCE: Institute of Nanoscience and Engineering, Department of Chemistry, University of Massachusetts Lowell, Lowell, MA, 01854, USA

SOURCE: Chemical Communications (Cambridge, United Kingdom) (2002), (17), 1854-1855

CODEN: CHCOFS; ISSN: 1359-7345

PUBLISHER: Royal Society of Chemistry

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 138:106347

AB The first, highly two-photon active C60 derivative comprised of a A-sp<sup>3</sup>-D conjugate structure was synthesized showing effective two-photon absorption cross-sections ( $\sigma_2' = 196 + 10^{-48} \text{ cm}^4 \text{ sec}^{-1} \text{ mol}^{-1}$ ) in the nanosecond regime among the best values for diphenylaminofluorene-based AFX chromophores.

IT 486998-58-7P

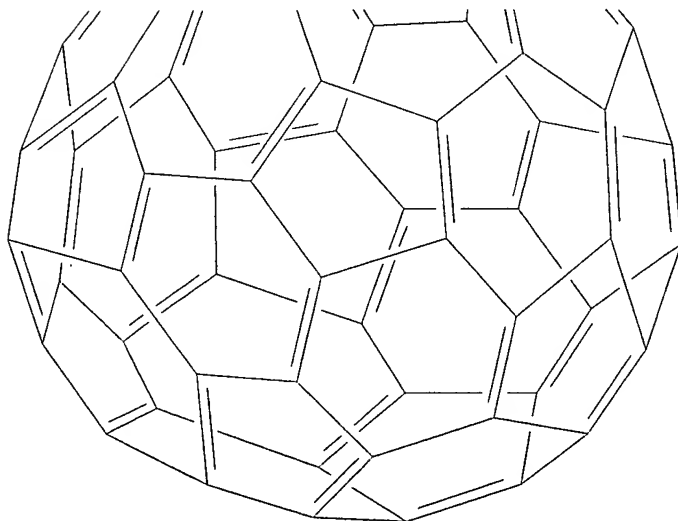
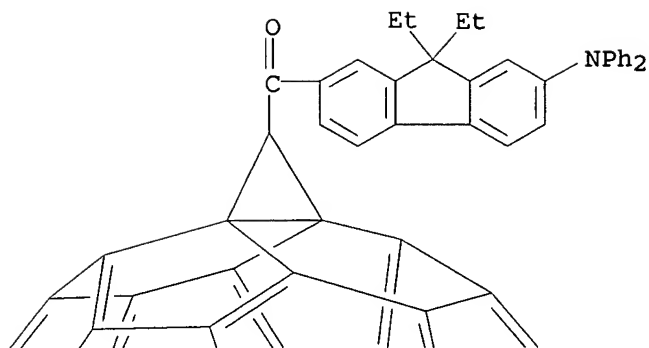
RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)

(2PA cross-section; synthesis of C60-diphenylaminofluorene dyad with large 2PA cross-sections and efficient intramol. two-photon energy transfer)

RN 486998-58-7 CAPLUS

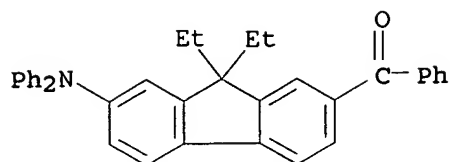
CN Methanone, 3'H-cyclopropa[1,9][5,6]fulleren-C60-1h-3'-yl[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)





IT 349123-45-1, AF-370  
 RL: PRP (Properties)  
 (2PA cross-section; synthesis of C60-diphenylaminofluorene dyad with  
 large 2PA cross-sections and efficient intramol. two-photon energy  
 transfer)  
 RN 349123-45-1 CAPLUS  
 CN Methanone, [7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]phenyl- (9CI)

(CA INDEX NAME)

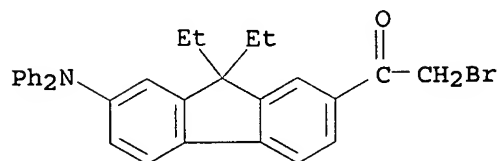


IT 486998-57-6P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (cyclopropanation; synthesis of C60-diphenylaminofluorene dyad with large 2PA cross-sections and efficient intramol. two-photon energy transfer)

RN 486998-57-6 CAPLUS

CN Ethanone, 2-bromo-1-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]- (9CI)  
 (CA INDEX NAME)



REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 43 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2002:458216 CAPLUS

DOCUMENT NUMBER: 137:39382

TITLE: 2,2'-Bridged biphenyl compound, optical recording material, and recording method

INVENTOR(S): Akiba, Masaharu

PATENT ASSIGNEE(S): Fuji Photo Film Co., Ltd., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 20 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent

LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2002172864	A2	20020618	JP 2001-110119	20010409
PRIORITY APPLN. INFO.:			JP 2000-297219	A 20000928
OTHER SOURCE(S):	MARPAT 137:39382			

GI

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB Laser-sensitive optical recording material contains I [X = CR1R2, NR3, O,

S, Se; R1-3 = alkyl, alkenyl, alkynyl, aryl, heterocycle (these may be substituted); R = substituent; n1-2 = 0-4]. The material is recorded by using  $\geq 2$  photon absorption induced by irradiation of laser beam with wavelength having no linear absorption and longer than linear absorption band of I. Five other 2,2'-bridged biphenyl compds. such as II, III, and IV [R1-2, R4, R13-14 = alkyl, alkenyl, alkynyl, aryl, heterocycle (these may be substituted); R5-8 = substituent; R15 = (substituted) heterocycle; XY- = Y-valent anion; Y = 1-5; h = 0-4; j = 0-6; i = 0-10; p = 0-2m; m = 0-5] are also claimed. The material shows high d. and capacity recording without using short-wavelength laser.

IT 437713-15-0P

RL: DEV (Device component use); PNU (Preparation, unclassified); PREP (Preparation); USES (Uses)  
(optical recording material containing bridged biphenyl compound)

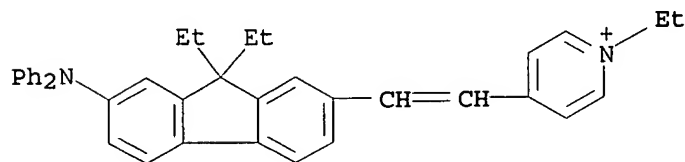
RN 437713-15-0 CAPLUS

CN Pyridinium, 4-[2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]-1-ethyl-, perchlorate (9CI) (CA INDEX NAME)

CM 1

CRN 437713-14-9

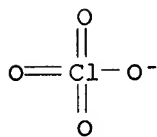
CMF C38 H37 N2



CM 2

CRN 14797-73-0

CMF Cl O4

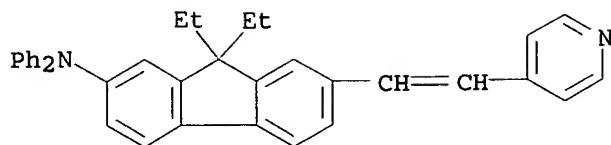


IT 197969-56-5

RL: DEV (Device component use); RCT (Reactant); RACT (Reactant or reagent); USES (Uses)  
(optical recording material containing bridged biphenyl compound)

RN 197969-56-5 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



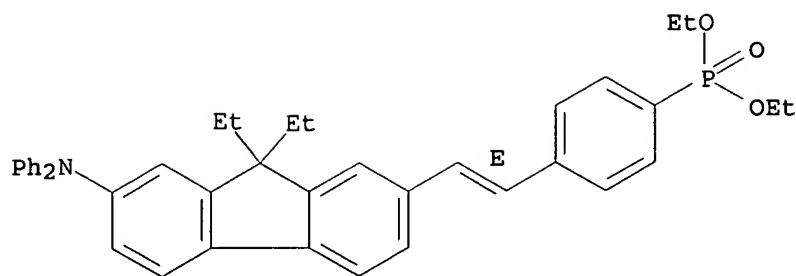
L6 ANSWER 44 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2002:372749 CAPLUS  
 DOCUMENT NUMBER: 137:176417  
 TITLE: Spectral properties of several fluorene derivatives with potential as two-photon fluorescent dyes  
 AUTHOR(S): Belfield, K. D.; Bondar, M. V.; Przhonska, O. V.; Schafer, K. J.; Mourad, W.  
 CORPORATE SOURCE: Department of Chemistry and CREOL/School of Optics, University of Central Florida, Orlando, FL, 32816-2366, USA  
 SOURCE: Journal of Luminescence (2002), 97(2), 141-146  
 CODEN: JLUMA8; ISSN: 0022-2313  
 PUBLISHER: Elsevier Science B.V.  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English

AB Investigations of the absorption, steady-state fluorescence, excitation and excitation anisotropy properties of several fluorene derivs., (7-benzothiazol-2-yl-9,9-didecylfluoren-2-yl)-diphenylamine, 9,9-didecyl-2,7-bis-(N,N-diphenylamino)fluorene and {4-[2-(7-diphenylamino-9,9-diethylfluoren-2-yl)vinyl]phenyl}phosphoric acid di-Et ester, in liquid solns. have been conducted. Spectral characteristics of these compds., including fluorescence quantum yields, were measured in acetonitrile, methylene chloride, THF and hexane at room temperature. Excitation anisotropy spectra provided a means to determine the nature of the short wavelength absorption bands as an electronic transition into a higher excited singlet state. It was found that excitation spectra in the short wavelength region do not correspond to the absorption bands that are correlated with the wavelength dependence of the fluorescence quantum yields. Major reasons of such spectral behavior are discussed.

IT 252000-05-8  
 RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); TEM (Technical or engineered material use); PROC (Process); USES (Uses)  
 (UV-visible absorption, fluorescence, excitation anisotropy and excitation spectra of fluorene derivs. with potential as two-photon fluorescent dyes in solvents of varying polarity)

RN 252000-05-8 CAPLUS  
 CN Phosphonic acid, [4-[(1E)-2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



REFERENCE COUNT: 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 45 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2001:923862 CAPLUS  
 DOCUMENT NUMBER: 136:54238  
 TITLE: Multiphoton photosensitization system  
 INVENTOR(S): Devoe, Robert J.  
 PATENT ASSIGNEE(S): 3M Innovative Properties Company, USA  
 SOURCE: PCT Int. Appl., 66 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

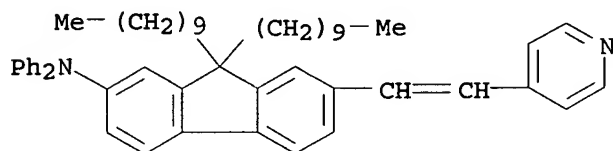
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001096409	A2	20011220	WO 2001-US19164	20010614
WO 2001096409	A3	20020404		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
EP 1297021	A2	20030402	EP 2001-946384	20010614
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
JP 2004503616	T2	20040205	JP 2002-510544	20010614
US 6852766	B1	20050208	US 2002-311041	20021212
US 2005054744	A1	20050310	US 2004-964510	20041013
PRIORITY APPLN. INFO.:			US 2000-211703P	P 20000615
			WO 2001-US19164	W 20010614
			US 2002-311041	A3 20021212

OTHER SOURCE(S): MARPAT 136:54238

AB A method of multiphoton photosensitizing a photoreactive composition comprises irradiating the composition with light sufficient to cause simultaneous absorption of at least two photons, thereby inducing at least one acid- or radical-initiated chemical reaction where the composition is exposed to the light.

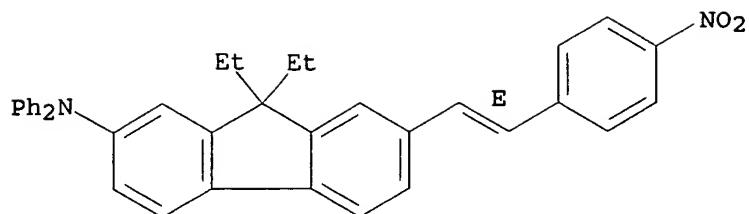
The composition comprises: (a) at least one reactive species that is capable of undergoing such reaction; and (b) at least one multi-component, multiphoton photoinitiator system.

IT 191667-13-7  
 RL: CAT (Catalyst use); USES (Uses)  
 (multiphoton photosensitization system)  
 RN 191667-13-7 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



L6 ANSWER 46 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2001:391256 CAPLUS  
 DOCUMENT NUMBER: 135:217791  
 TITLE: Effects of  $\pi$  centers and symmetry on two-photon  
 absorption cross sections of organic chromophores  
 AUTHOR(S): Wang, Chuan-Kui; Macak, Peter; Luo, Yi; Agren, Hans  
 CORPORATE SOURCE: Theoretical Chemistry, Royal Institute of Technology,  
 Stockholm, S-10044, Swed.  
 SOURCE: Journal of Chemical Physics (2001), 114(22), 9813-9820  
 CODEN: JCPSA6; ISSN: 0021-9606  
 PUBLISHER: American Institute of Physics  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 AB The authors have theor. examined organic mols. that exhibit large 2-photon  
 absorption cross sections in the visible region and that were synthesized  
 in different labs. One- and 2-photon absorption cross sections of the 4  
 lowest excited states of each mol. were calculated at the same theor. level  
 using ab initio response theory. The mol. length and the 1-photon  
 absorption intensity are quite strongly correlated factors, but a  
 corresponding correlation for the 2-photon absorption is much weaker or is  
 missing. But a most crucial role for large 2-photon absorption is played  
 by the  $\pi$  center. For mols. with a given  $\pi$  center a sym. structure  
 with strong donor groups can result in a maximum 2-photon absorption cross  
 section. Theor. findings are consistent with some recent exptl.  
 observations. The chromophore based on dithienothiophene as  $\pi$  center  
 attached with sym. N,N-diphenylamine donors has the largest 2-photon cross  
 section in the visible region among all known 1-dimensional 2-photon organic  
 materials that were reported in the literature.  
 IT 252000-06-9 358374-63-7  
 RL: PRP (Properties)  
 (effects of  $\pi$  centers and symmetry on two-photon absorption cross  
 sections of organic chromophores)  
 RN 252000-06-9 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-7-[(1E)-2-(4-nitrophenyl)ethenyl]-N,N-  
 diphenyl- (9CI) (CA INDEX NAME)

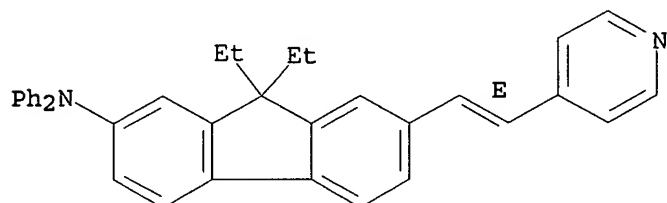
Double bond geometry as shown.



RN 358374-63-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[(1E)-2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 47 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2001:298453 CAPLUS

DOCUMENT NUMBER: 135:93912

TITLE: Diphenylaminofluorene-Based Two-Photon-Absorbing Chromophores with Various  $\pi$ -Electron Acceptors

AUTHOR(S): Kannan, Ramamurthi; He, Guang S.; Yuan, Lixiang; Xu, Faming; Prasad, Paras N.; Dombroskie, Ann G.; Reinhardt, Bruce A.; Baur, Jeffery W.; Vaia, Richard A.; Tan, Loon-Seng

CORPORATE SOURCE: Systran Systems Corporation, Dayton, OH, 45432, USA

SOURCE: Chemistry of Materials (2001), 13(5), 1896-1904

CODEN: CMATEX; ISSN: 0897-4756

PUBLISHER: American Chemical Society

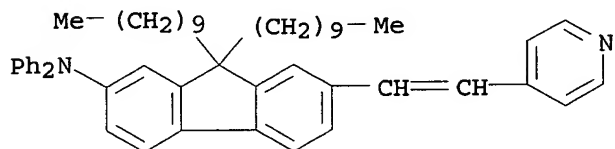
DOCUMENT TYPE: Journal

LANGUAGE: English

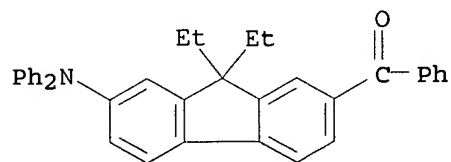
AB A new series of linear, asym. (diphenylamino)fluorene-based chromophores (AFX) with various strong  $\pi$ -electron acceptors were synthesized and evaluated for two-photon absorptivity. These chromophores were studied to determine a suitable replacement for 2-(4-pyridinyl)vinyl, the  $\pi$  acceptor for our previously reported AFX series, which contains a photochem. and thermooxidatively unstable olefinic unit. In addition to the benzoyl group (AF-370), these  $\pi$ -electron acceptors include 2-benzothiazolyl (AF-240), 2-benzoxazolyl (AF-390), N-phenyl-2-benzimidazolyl (AF-386), and 3,4-diphenyl-1H-imidazol-2-yl (AF-385) moieties (five-membered heterocycles) and the 2-quinoxaliny (AF-260) group (six-membered heterocycle). From nanosecond nonlinear transmission measurements, these new chromophores have effective two-photon cross sections ( $\sigma_2'$ ) at 800 nm spanning from  $3.87 \times 10^{-48} \text{ cm}^4 \text{ s}/(\text{photon mol.})$  for AF-385 to  $97.46 \times 10^{-48} \text{ cm}^4 \text{ s}/(\text{photon mol.})$  for AF-240. Two of them, AF-240 and AF-370 [ $\sigma_2' = 84.32 \times 10^{-48} \text{ cm}^4 \text{ s}/(\text{photon mol.})$ ], stand out as having relatively good, albeit lower, values of two-photon cross

sections, as compared to that of previously reported N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-didecyl-2-fluorenamine (AF-50) [ $\sigma_2' = 115.6 + 10.48 \text{ cm}^4 \text{ s}/(\text{photon mol.})$ ]. However, we observed that AF-240 was more photochem. robust than AF-50 when their THF solns. were subjected to repetitive and prolonged exposure to nanosecond laser radiation. On the basis of nanosecond TPA cross-section data ( $\sigma_2'/\text{mol. weight values}$ ), the general trend for  $\pi$ -electron accepting ability, i.e., ability to accept charge transferred from diphenylamine, appears to be as follows: 2-(4-pyridinyl)vinyl > 2-benzothiazolyl > benzoyl > N-phenyl-2-benzimidazolyl > 2-quinoxalinylyl > 2-benzoxazolyl > 4,5-diphenyl-2-imidazolyl.

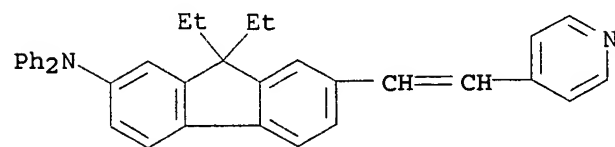
IT 191667-13-7, AF 50  
 RL: PRP (Properties)  
 ((diphenylamino)fluorene-based two-photon-absorbing chromophores with various  $\pi$ -electron acceptors)  
 RN 191667-13-7 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



IT 349123-45-1P  
 RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
 (AF 370; preparation of (diphenylamino)fluorene-based two-photon-absorbing chromophores with various  $\pi$ -electron acceptors)  
 RN 349123-45-1 CAPLUS  
 CN Methanone, [7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]phenyl- (9CI) (CA INDEX NAME)

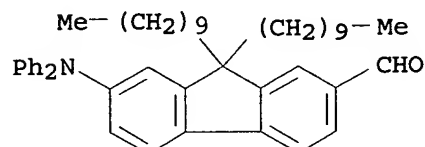


IT 197969-56-5  
 RL: PRP (Properties)  
 (AF 60; (diphenylamino)fluorene-based two-photon-absorbing chromophores with various  $\pi$ -electron acceptors)  
 RN 197969-56-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)





IT 348631-00-5P  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT  
 (Reactant or reagent)  
 (preparation of (diphenylamino)fluorene-based two-photon-absorbing  
 chromophores with various  $\pi$ -electron acceptors)  
 RN 348631-00-5 CAPLUS  
 CN 9H-Fluorene-2-carboxaldehyde, 9,9-didecyl-7-(diphenylamino)- (9CI) (CA  
 INDEX NAME)

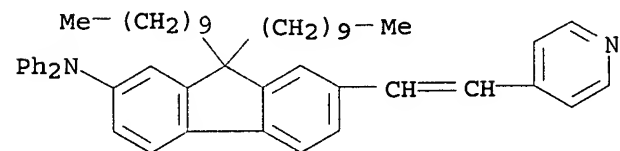


REFERENCE COUNT: 54 THERE ARE 54 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 48 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2000:791214 CAPLUS  
 DOCUMENT NUMBER: 133:355509  
 TITLE: Theoretical model for excited-state absorption  
 AUTHOR(S): Das, G. P.; Yeates, A. T.; Dudis, D. S.  
 CORPORATE SOURCE: U.S. Air Force Research Lab, MLBP Wright Peterson AFB,  
 OH, 45433, USA  
 SOURCE: International Journal of Quantum Chemistry (2000),  
 80(4/5), 1039-1042  
 CODEN: IJQCB2; ISSN: 0020-7608  
 PUBLISHER: John Wiley & Sons, Inc.  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English

AB For long incident pulses, measurement of two-photon absorption coeffs. can  
 be complicated by excited-state absorption. We develop here a quant. ab  
 initio model to determine the contribution of excited-state absorption in the  
 interpretation of the two-photon absorption data for such materials. We  
 present representative calcns. of this effect for some selected systems.

IT 191667-13-7, AF50  
 RL: PEP (Physical, engineering or chemical process); PROC (Process)  
 (theor. model for excited-state absorption)  
 RN 191667-13-7 CAPLUS  
 CN 9H-Fluorene-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 49 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 2000:766204 CAPLUS  
 DOCUMENT NUMBER: 134:107574  
 TITLE: Multiphoton absorption and optical limiting

AUTHOR(S): Sutherland, R. L.; McLean, D. G.; Kirkpatrick, S. M.;  
Fleitz, P. A.; Chandra, S.; Brant, M. C.

CORPORATE SOURCE: Air Force Research Laboratory, Materials and  
Manufacturing Directorate, Wright-Patterson Air Force  
Base, OH, 45433-7750, USA

SOURCE: NATO Science Series, 3: High Technology (2000),  
79(Multiphoton and Light Driven Multielectron  
Processes in Organics), 67-81  
CODEN: NSSTFF; ISSN: 1388-6576

PUBLISHER: Kluwer Academic Publishers

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Optical limiting requires the development of materials exhibiting strong  
nonlinear absorption. Multiphoton absorption in materials can take the  
form of instantaneous two-photon absorption or time-integrating excited  
state absorption. The authors examining both types of mechanisms for optical  
limiting. Several exptl. methods were used to study these phenomena and  
characterize materials. Often new or unusual effects arise in these  
expts. that require careful anal. to elucidate desired parameters. The  
authors describe several of these and give results for different materials  
relevant to optical limiting.

IT 191667-13-7, AF-50

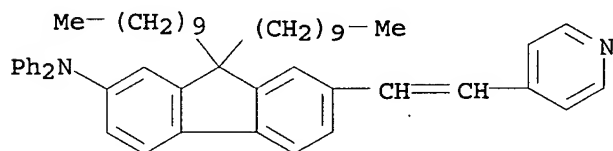
RL: PEP (Physical, engineering or chemical process); PRP (Properties);

PROC (Process)

(multiphoton absorption and optical limiting and nonlinear optical  
properties for several materials)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS  
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 50 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:606926 CAPLUS

DOCUMENT NUMBER: 133:200634

TITLE: Nonlinear spectrophotometer

INVENTOR(S): Mukherjee, Anadi; Mukherjee, Nandini; Sarvis, Conrad  
S.; Reinhardt, Bruce A.

PATENT ASSIGNEE(S): United States Dept. of the Air Force, USA

SOURCE: U.S., 9 pp.

CODEN: USXXAM

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

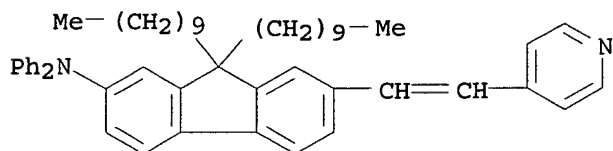
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6111641	A	20000829	US 1998-46462	19980323
US 5991021	A	19991123	US 1999-241191	19990201
PRIORITY APPLN. INFO.:			US 1998-46462	A3 19980323

AB Nonlinear spectrophotometers for fluorescent materials are described which comprise a laser for producing a primary laser beam; a second harmonic generation crystal that alters the primary laser beam; a beam splitter for dividing the primary laser beam into a first signal manifold laser beam and a second signal manifold laser beam; a first neutral d. filter for filtering the first signal manifold laser beam to a first harmonic spectrum; a means for passing filtered first signal manifold laser beam through a sample; a first sensor for detecting filtered first signal manifold laser beam after passing through a sample; a second neutral d. filter for filtering the second signal manifold laser beam to a second harmonic spectrum; a means for passing filtered second signal manifold laser beam through a sample; a second sensor for detecting filtered second signal manifold laser beam after passing through a sample; and a means to compare the outputs of the first and second sensors. The spectrophotometers may be used to measure two-photon absorption coeffs. and cross-sections.

IT 191667-13-7, AF-50  
 RL: ANT (Analyte); PRP (Properties); ANST (Analytical study)  
 (nonlinear spectrophotometers)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



L6 ANSWER 51 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:463667 CAPLUS

DOCUMENT NUMBER: 133:184925

TITLE: Photoluminescence studies and read/write process of a strong two-photon absorbing chromophore

AUTHOR(S): Sivaraman, R.; Clarson, S. J.; Lee, B. K.; Steckl, A. J.; Reinhardt, B. A.

CORPORATE SOURCE: Department of Materials Science and Engineering, University of Cincinnati, Cincinnati, OH, 45221-0012, USA

SOURCE: Applied Physics Letters (2000), 77(3), 328-330  
 CODEN: APPLAB; ISSN: 0003-6951

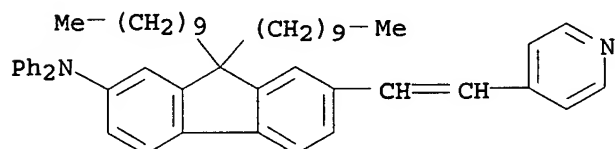
PUBLISHER: American Institute of Physics

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The optical properties of a fluorene-based chromophore exhibiting a high two-photon absorption cross section were studied both in the pure state and as a guest-host system in poly(N-vinylcarbazole). Upon irradiation with a He Cd (He-Cd) laser at 325 nm, the guest-host sample exhibits a blueshift with a maximum emission at 459 nm. Information written onto the samples in the blue can also be read using an Ar+ laser. The chromophore undergoes a chemical change upon irradiation in air, and was found to no longer exhibit upconversion after this transformation. IR anal. of the sample (the chromophore) before and after irradiation was carried out, and the spectra suggest the formation of a new conjugated species. Other applications the authors have developed using this phenomenon include the successful writing of optical device structures in polymer films where the AF-50 is a guest material.

IT 191667-13-7  
 RL: PEP (Physical, engineering or chemical process); PRP (Properties);  
 PROC (Process)  
 (photoluminescence studies and read/write process of a strong  
 two-photon absorbing chromophore)  
 RN 191667-13-7 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 52 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1999:748391 CAPLUS  
 DOCUMENT NUMBER: 131:344066  
 TITLE: Nonlinear spectrophotometer  
 INVENTOR(S): Mukherjee, Anadi; Mukherjee, Nandini; Sarvis, Conrad  
 S.; Reinhardt, Bruce A.  
 PATENT ASSIGNEE(S): United States Dept. of the Air Force, USA  
 SOURCE: U.S., 9 pp., Division of U.S. Ser. No. 46,462.  
 CODEN: USXXAM  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 2  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5991021	A	19991123	US 1999-241191	19990201
US 6111641	A	20000829	US 1998-46462	19980323
PRIORITY APPLN. INFO.:			US 1998-46462	A3 19980323

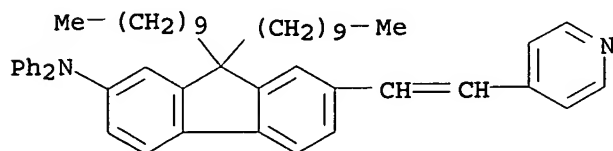
AB Combined nonlinear spectrophotometers for measuring both fluorescent and nonfluorescent materials are described which comprise a high intensity laser pump source; means for attenuating the the pump beam; means for dividing the pump beam into a first compartment pump beam and a second compartment pump beam; a first compartment nonlinear spectrophotometer for fluorescent sample measurements; and a second compartment nonlinear Michelson interferometer, energized by the second compartment pump beam, for measuring nonfluorescent samples and outputting a signal; means for receiving user inputs on sample parameters; and means for receiving signals from the first and second compartments and processing results. The first compartment nonlinear spectrophotometer includes a frequency doubling crystal that alters the first compartment pump beam; a beam splitter for dividing the first compartment pump beam into a first signal manifold laser beam and a second signal manifold laser beam; first and second neutral d. filters for filtering, resp., the first and second signal manifold laser beam to first and second harmonic spectrum; means for passing the filtered signal manifold beams through a sample; and sensors for detecting the filtered signal manifold beams after they pass through a sample. The instruments can be used to measure the two-photon absorption coeffs. and cross-sections of materials.

IT 191667-13-7, AF-50

RL: ANT (Analyte); PRP (Properties); ANST (Analytical study)  
(combined nonlinear spectrophotometers for measuring both fluorescent and nonfluorescent materials)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 53 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1999:634693 CAPLUS

DOCUMENT NUMBER: 132:28209

TITLE: New Two-Photon Absorbing Fluorene Derivatives:  
Synthesis and Nonlinear Optical Characterization  
AUTHOR(S): Belfield, Kevin D.; Hagan, David J.; Van Stryland,  
Eric W.; Schafer, Katherine J.; Negres, Raluca A.

CORPORATE SOURCE: Department of Chemistry and School of Optics,  
University of Central Florida, Orlando, FL,  
32816-2366, USA

SOURCE: Organic Letters (1999), 1(10), 1575-1578  
CODEN: ORLEF7; ISSN: 1523-7060

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Efficient Pd-catalyzed Heck coupling methodol. was employed to provide 2 new fluorene derivs. with phosphonate (2) and nitro (3) electron-withdrawing functionalities. Both derivs. exhibit 2-photon absorption (2PA), as determined by nonlinear absorption measurements using a femtosecond pump/white light continuum probe NLO spectrometer. Both fluorene derivs. have high 2PA cross sections (650 and 1300 + 10-50 cm4 s. photon-1 mol.-1 for compds. 2 and 3, resp.).

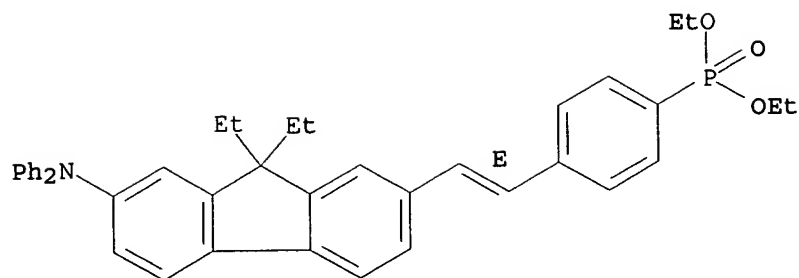
IT 252000-05-8P 252000-06-9P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
(new two-photon absorbing fluorene derivs. with synthesis, nonlinear optical absorptivity, and UV-visible spectra)

RN 252000-05-8 CAPLUS

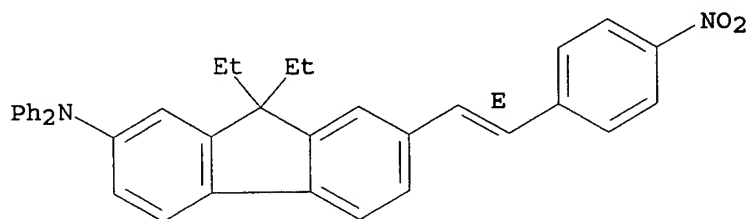
CN Phosphonic acid, [4-[(1E)-2-[7-(diphenylamino)-9,9-diethyl-9H-fluoren-2-yl]ethenyl]phenyl]-, diethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 252000-06-9 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-7-[(1E)-2-(4-nitrophenyl)ethenyl]-N,N-diphenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



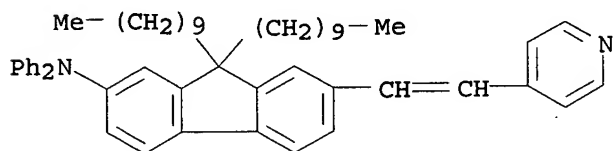
REFERENCE COUNT: 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 54 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1999:611786 CAPLUS  
 DOCUMENT NUMBER: 131:329489  
 TITLE: Molecular Environment Effects on Two-Photon-Absorbing Heterocyclic Chromophores  
 AUTHOR(S): Baur, Jeffery W.; Alexander, Max D. Jr.; Banach, Michael; Denny, Lisa R.; Reinhardt, Bruce A.; Vaia, Richard A.; Fleitz, Paul A.; Kirkpatrick, Sean M.  
 CORPORATE SOURCE: Polymer Branch AFRL/MLBP, U.S. Air Force Research Laboratory, Wright-Patterson AFB, OH, 45433, USA  
 SOURCE: Chemistry of Materials (1999), 11(10), 2899-2906  
 CODEN: CMATEX; ISSN: 0897-4756  
 PUBLISHER: American Chemical Society  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English

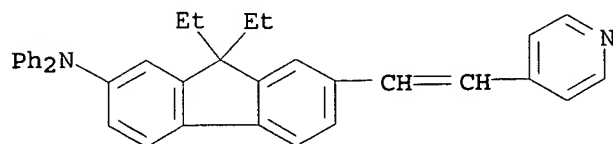
AB Over the past several years, organic mols. exhibiting significant 2-photon absorbance and subsequent up-converted fluorescence were of intense interest for a wide variety of applications including data storage, imaging, and optical limiting. The establishment of structure-property relations for some asym. mols. was hindered by the sensitivity of these nonlinear optical properties to the local mol. environment and to the pulse width of the incident radiation. To understand the influence of the local mol. environment on the excited states of these 2-photon-absorbing mols., the linear absorbance, the single-photon-excited luminescence, and the 2-photon-excited luminescence of heterocyclic dyes are examined. The stabilization of the longest-lived 1-photon-excited state by the local mol. environment can be described by mean field interactions with solvent mols. as given by the Lippert equation. Because the same stabilization

dominates the 2-photon-induced longest-lived excited state, the influence of the local mol. environment on the 2-photon luminescence can be predicted using the Lippert equation and 1-photon expts. These results support models that suggest excited-state absorption is the primary cause of sensitivity of the effective 2-photon cross-section to the pulse-width and the local mol. environment.

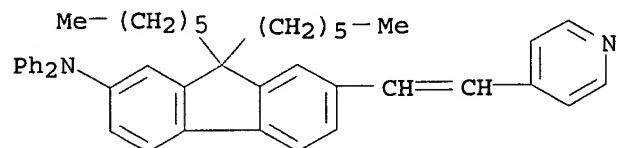
IT 191667-13-7, N,N-Diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-didecylfluoren-2-amine 197969-56-5, N,N-Diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-diethylfluoren-2-amine 209603-50-9, N,N-Diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-dihexylfluoren-2-amine  
 RL: PEP (Physical, engineering or chemical process); PRP (Properties); PROC (Process)  
 (mol. environment effects on two-photon-absorbing)  
 RN 191667-13-7 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



RN 197969-56-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



RN 209603-50-9 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-dihexyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 25 THERE ARE 25 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 55 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1999:532016 CAPLUS  
 DOCUMENT NUMBER: 131:206638  
 TITLE: Measurement of two-photon absorption coefficients employing femtosecond, phase-mismatched DFWM, and nDFWM spectroscopy  
 AUTHOR(S): Dalton, Larry R.; Larsen, Rulon J.; Strohkendl, Fritz  
 CORPORATE SOURCE: Loker Hydrocarbon Research Institute, Univ. Southern

SOURCE: California, Los Angeles, CA, USA  
MCLC S&T, Section B: Nonlinear Optics (1999), 21(1-4),  
245-261  
CODEN: MCLOEB; ISSN: 1058-7268

PUBLISHER: Gordon & Breach Science Publishers

DOCUMENT TYPE: Journal

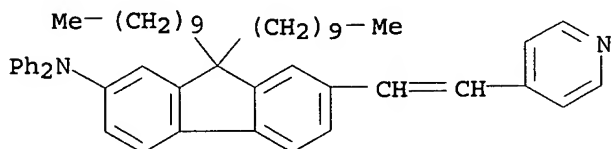
LANGUAGE: English

AB The real and imaginary components of the 3rd order nonlinear optical susceptibility for the fullerenes (C60, C70) and for N,N-diphenyl-7-[2-(4-pyridinyl) ethenyl]-9,9-di-n-decyl-fluoren-2-amine (AF50) were measured over the wavelength range from 0.5-2.0  $\mu\text{m}$  by a variety of techniques (including newly developed techniques of nDFWM and THG/DFWM) based on 4-wave mixing employing 100 fs pulses. Results are compared with related data obtained by 3rd harmonic generation, electroabsorption, and direct 2-photon absorption measurements. The magnitudes of 2-photon absorption for these materials are found to be comparable to that of GaAs (0.020 cm/MW).

IT 191667-13-7, AF50  
RL: PRP (Properties)  
(two-photon absorption coeffs. of AF50 employing femtosecond, phase-mismatched DFWM, and nDFWM spectroscopy)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



REFERENCE COUNT: 18 THERE ARE 18 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 56 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1999:107637 CAPLUS

DOCUMENT NUMBER: 130:259118

TITLE: Femtosecond, Frequency-Agile, Phase-Sensitive-Detected, Multi-Wave-Mixing Nonlinear Optical Spectroscopy Applied to  $\pi$ -Electron Photonic Materials

AUTHOR(S): Drenser, K. A.; Larsen, R. J.; Strohkendl, F. P.; Dalton, L. R.

CORPORATE SOURCE: Loker Hydrocarbon Research Institute, University of Southern California, CA, 90089-1662, USA

SOURCE: Journal of Physical Chemistry A (1999), 103(14), 2290-2301  
CODEN: JPACFH; ISSN: 1089-5639

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Degenerate four-wave mixing (DFWM) spectroscopy is modified to exploit femtosecond pulses, phase-sensitive-detection, frequency (wavelength) agility, two-color (nearly degenerate multi-wave mixing) radiation, and improved signal-to-noise capabilities that can be realized through a combination of new solid state lasers, nonlinear optical components, and novel design concepts. The resulting time-resolved nonlinear optical techniques permit instantaneous optical nonlinearities, such as two-photon



absorption cross sections, to be accurately measured over the spectral range from 450 to 2500 nm (and with significantly greater effort from 225 to 5000 nm). The power of the new techniques is illustrated by their application to the definition of Hg two-photon resonances of C60 and C70 as well as to the characterization of optical nonlinearities in two linear chromophores of putative utility for sensor protection and electrooptic modulation. Explicitly, these measurements provide accurate determination of

both

transition energies and transition moments (matrix elements connecting the two photon levels). Results are compared to those previously reported in the literature illustrating the advantages and problems associated with particular measurement techniques. All of the mols. studied exhibit two-photon absorption coeffs. comparable to that of GaAs, the most studied putative sensor protection material (based on use of electronic optical nonlinearity). Femtosecond pulse techniques are shown, in all cases, to be necessary to avoid complications arising from excited-state absorption and relaxation phenomena. The importance of phase-sensitive detection in identifying complications from overlapping transitions is illustrated.

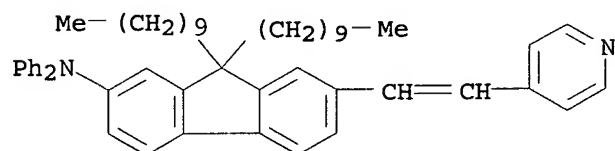
IT 191667-13-7

RL: PRP (Properties)

(femtosecond, frequency-agile, phase-sensitive-detected,  
multi-wave-mixing nonlinear optical spectroscopy applied to  
 $\pi$ -electron photonic materials)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



REFERENCE COUNT: 114 THERE ARE 114 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 57 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:805856 CAPLUS

DOCUMENT NUMBER: 130:145769

TITLE: Nonlinear measurements on AF-50

AUTHOR(S): Fleitz, P. A.; Brant, M. C.; Sutherland, R. L.;  
Strohkendl, F. P.; Larsen, R. J.; Dalton, L. R.

CORPORATE SOURCE: Materials Directorate, Air Force Research Laboratory,  
AFRL/MLPJ, Wright Patterson Air Force Base, OH,  
45433-7702, USA

SOURCE: Proceedings of SPIE-The International Society for  
Optical Engineering (1998), 3472(Nonlinear Optical  
Liquids for Power Limiting and Imaging), 91-97  
CODEN: PSISDG; ISSN: 0277-786X

PUBLISHER: SPIE-The International Society for Optical Engineering

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The nonlinear optical properties of N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-di-n-decylfluoren-2-amine (AF-50) were investigated. The nonlinear absorption of a saturated solution of this material in Me2CO was investigated with 430 fs pulses at 790 nm. From these results, the 2-photon absorption cross-section was determined to be  $25 \pm 10^{-50}$

cm<sup>4</sup>sec/photon mol. This number is in agreement (within a factor of 2) with theor. calcns. Nonlinear absorption and optical limiting measurements were also made using a Nd:YAG pumped dye laser with 4.3 ns pulses at 694 nm. These results suggest inherent differences in the performance of 2-photon absorbing materials in these 2 different geometries.

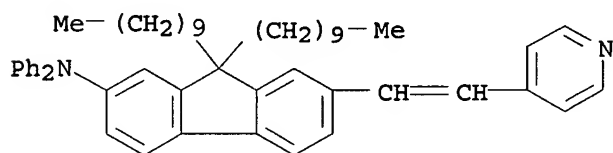
IT 191667-13-7, AF-50

RL: PEP (Physical, engineering or chemical process); PRP (Properties); PROC (Process)

(nonlinear optical absorption measurement and 2-photon absorption cross-section calcn. of AF-50)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



REFERENCE COUNT: 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 58 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:805854 CAPLUS

DOCUMENT NUMBER: 130:139023

TITLE: Influence of molecular environment on single photon behavior of heterocyclic NLO chromophores and its implications to two-photon behavior

AUTHOR(S): Baur, Jeffery W.; Alexander, Max D., Jr.; Banach, Michael; Reinhardt, Bruce; Prasad, Paras N.; Yaun, Lixiang; Vaia, Richard A.

CORPORATE SOURCE: Materials and Manufacturing Directorate, Air Force Research Laboratory, Wright-Patterson AFB, OH, 45433, USA

SOURCE: Proceedings of SPIE-The International Society for Optical Engineering (1998), 3472(Nonlinear Optical Liquids for Power Limiting and Imaging), 70-79  
CODEN: PSISDG; ISSN: 0277-786X

PUBLISHER: SPIE-The International Society for Optical Engineering

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The influence of the solvent environment on linear absorbance and photoluminescence of a series of donor-acceptor heterocyclic chromophores (AF50, AF100, AF60, AF210) was examined. The Stoke's shift associated with 1-photon absorbance and photoluminescence was observed to increase with increasing solvent polarity. This behavior is adequately described by the Lippert equation and is related to relaxation of the solvent mols. around an excited mol. Addnl., it was observed that the spectral shape, as well as the solvent dependence, of 2-photon and 1-photon pumped photoluminescence were similar, thus indicating that the longest-lived luminescing excited state is independent of the method of excitation. These results have direct implications to 2-photon applications which rely on up-converted fluorescence. They also yield insight into the structure-property relationships governing their linear and multi-photon behavior including the potential contributions to the effective 2-photon cross-section from excited state absorbance.

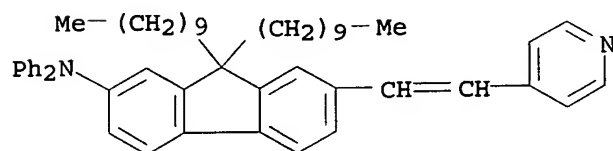
IT 191667-13-7, AF50 197969-56-5, 9H-Fluoren-2-amine,

9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- 209603-50-9  
 , 9H-Fluoren-2-amine, 9,9-dihexyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 RL: PEP (Physical, engineering or chemical process); PRP (Properties);  
 PROC (Process)

(mol. environment effect on single photon absorptivity and luminescence  
 of heterocyclic NLO chromophores and its implications to 2-photon  
 behavior)

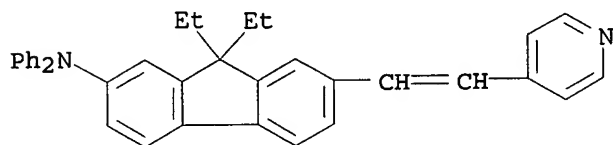
RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



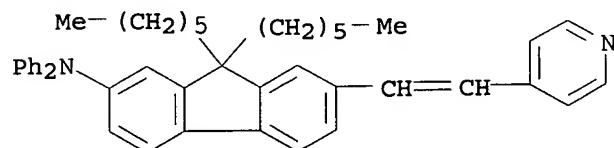
RN 197969-56-5 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



RN 209603-50-9 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-dihexyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 59 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:781324 CAPLUS

DOCUMENT NUMBER: 130:145737

TITLE: Probing two-photon excitation dynamics using ultrafast  
 laser pulses

AUTHOR(S): Swiatkiewicz, J.; Prasad, P. N.; Reinhardt, B. A.

CORPORATE SOURCE: Photonics Research Laboratory, Departments of  
 Chemistry and Physics, State University of New York,  
 Buffalo, NY, 14260-3000, USA

SOURCE: Optics Communications (1998), 157(1-6), 135-138

CODEN: OPCOB8; ISSN: 0030-4018

PUBLISHER: Elsevier Science B.V.

DOCUMENT TYPE: Journal

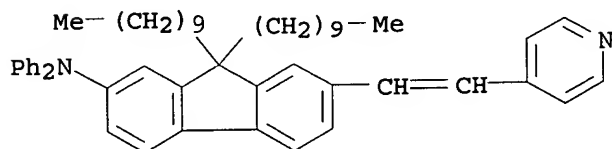
LANGUAGE: English

AB The authors probe the two-photon excitation dynamics of two new dyes, N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-di-n-decyl-fluoren-2-amine (AF50) and (7-(7-benzothiazol-2-yl-9,9-diethylfluoren-2-yl)-9,9-diethylfluoren-2-yl)diphenylamine (AF250) using femtosecond excitation pulses by Z-scan and time-resolved pump-probe absorption measurements. Irradiance dependence of the induced absorption cross-section is linked to linear absorption of the two-photon excited state. The excited state linear absorption cross-section are  $1.0 \times 10^{-17} \text{ cm}^2$  for AF250 and  $2.7 \times 10^{-17} \text{ cm}^2$  for AF250. Relaxation of the two-photon excited state follows a complicated path with three distinct relaxation times. The longest ones, 1.6 ns for the AF50 and 1.9 ns for the AF250, are associated with the resp. lowest singlet life-times: 2.23 ns and 2.15 ns.

IT 191667-13-7  
RL: PRP (Properties)  
(probing two-photon excitation dynamics using ultrafast laser pulses)

RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 60 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:532424 CAPLUS

DOCUMENT NUMBER: 129:308216

TITLE: Two-photon processes: dynamics and applications to photonics and biophotonics

AUTHOR(S): Prasad, P. N.; Joshi, M.; Haridas, E. P.; Swiatkiewicz, J.; He, G. S.; Bhawalkar, J. D.; Kim, K. S.; Reinhardt, B. A.

CORPORATE SOURCE: Photonics Res. Lab., Dep. Chem. and Phys., State Univ. New York, Buffalo, NY, 14260, USA

SOURCE: Polymer Preprints (American Chemical Society, Division of Polymer Chemistry) (1998), 39(2), 1114-1115  
CODEN: ACPPAY; ISSN: 0032-3934

PUBLISHER: American Chemical Society, Division of Polymer Chemistry

DOCUMENT TYPE: Journal

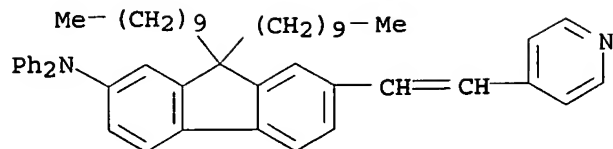
LANGUAGE: English

AB The authors have presented the authors' effort to synthesize mols. and bulk systems with very high two-photon properties. The authors have also carried out pump-probe expts. which reveal the dynamics of multiphoton absorption. Finally the authors demonstrated various applications of two-photon process that are being carried out in the authors' laboratory

IT 191667-13-7, AF 50  
RL: PRP (Properties)  
(two-photon processes for dye mols. and their applications to photonics and biophotonics)

RN 191667-13-7 CAPLUS

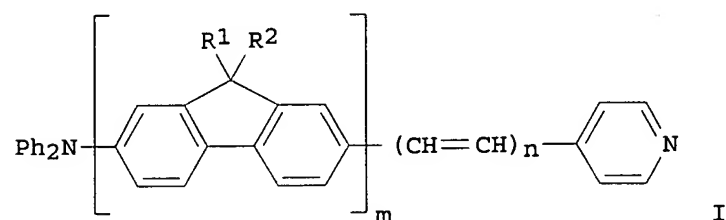
CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 61 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1998:427808 CAPLUS  
 DOCUMENT NUMBER: 129:96587  
 TITLE: Asymmetric dyes with large two-photon absorption cross sections  
 INVENTOR(S): Reinhardt, Bruce A.; Bhatt, Jayprakash C.; Brott, Lawrence L.; Clarson, Stephen J.  
 PATENT ASSIGNEE(S): United States Dept. of the Air Force, USA  
 SOURCE: U.S., 5 pp.  
 CODEN: USXXAM  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

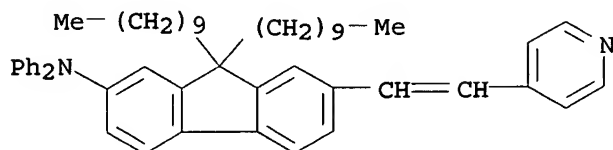
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5770737	A	19980623	US 1997-933067	19970918
PRIORITY APPLN. INFO.:			US 1997-933067	19970918
OTHER SOURCE(S):	MARPAT 129:96587			
GI				



AB A two-photon absorbing chromophore, useful in laser-scanning confocal fluorescence microscopy, has the structure I [R1, R2 = C8-12 alkyl; m = 1, 2; n = 0, 1]. Thus, fluorene was treated in 2 stages with BuLi followed by Me(CH2)9Br and the product was brominated in CH2Cl2 in the presence of I2 to give 2,7-dibromo-9,9-didecylfluorene (II). II was condensed first with Ph2NLi in the presence of (2-MeC6H4)3P and a Pd complex and then with 4-vinylpyridine in the presence of (2-MeC6H4)3P, Et3N, and Pd(OAc)2 to give I (R1 = R2 = decyl, m = n = 1), a 0.0450M solution of which in THF showed a 2-photon absorption cross section of 78 + 10-20 cm4/GW when laser-irradiated at 800 nm and pulse duration 8-10 ns.

IT 191667-13-7P  
 RL: PRP (Properties); SPN (Synthetic preparation); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)  
 (asym. dyes with large two-photon absorption cross sections)  
 RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS  
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 62 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:411176 CAPLUS

DOCUMENT NUMBER: 129:96579

TITLE: Highly active two-photon dyes: design, synthesis, and  
characterization toward application

AUTHOR(S): Reinhardt, Bruce A.; Brott, Lawrence L.; Clarson,  
Stephen J.; Dillard, Ann G.; Bhatt, Jayprakash C.;  
Kannan, Ramamurthi; Yuan, Lixiang; He, Guang S.;  
Prasad, Paras N.

CORPORATE SOURCE: Polymer Branch WL/MLBP Materials Directorate, U. S.  
Air Force Research Laboratory, Wright-Patterson AFB,  
OH, 45433-7750, USA

SOURCE: Chemistry of Materials (1998), 10(7), 1863-1874  
CODEN: CMATEX; ISSN: 0897-4756

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

AB A series of compds. with systematically varied mol. structures which  
exhibit very large effective two-photon cross sections has been  
synthesized and characterized in solution using a nonlinear transmission  
technique. The general structure of these compds. can be categorized into  
two basic structural families: acceptor/donor/donor/acceptor and  
donor/bridge/acceptor. This study attempts to determine certain mol.  
structure/effective two-photon absorption relationships by careful  
characterization and as a function of systematically varied changes in the  
organic structure of the dye mols. Such information can be useful in the  
design of more efficient two-photon dyes for imaging and power-limiting  
applications. The results of the study indicate that with the  
incorporation of certain combinations of structural elements, dyes can be  
synthesized which have greatly increased effective cross sections as high  
as  $152.5 \times 10^{-48} \text{ cm}^4 \text{ s/photon mol.}$  in benzene solution at 800 nm using  
8-ns pulses. This value is orders of magnitude higher than com. available  
organic dyes measured at the same wavelength. Although the process is  
thought to involve a combination of two-photon absorption and excited  
state absorption phenomena, the information gathered from these new  
families of dyes has provided an important first step in producing  
improved materials for use in many different two-photon technol.  
application.

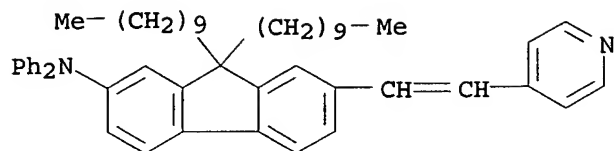
IT 191667-13-7P 197969-56-5P 209603-48-5P

209603-50-9P 209603-53-2P 209603-56-5P

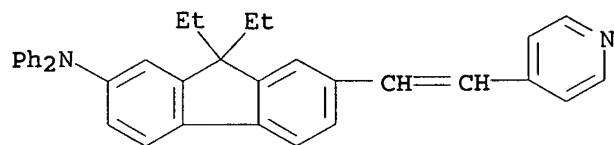
RL: PRP (Properties); SPN (Synthetic preparation); TEM (Technical or  
engineered material use); PREP (Preparation); USES (Uses)  
(preparation of highly active 2-photon dyes)

RN 191667-13-7 CAPLUS

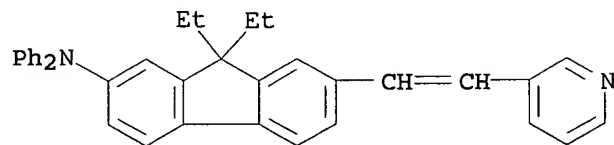
CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



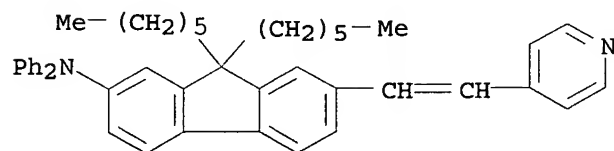
RN 197969-56-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



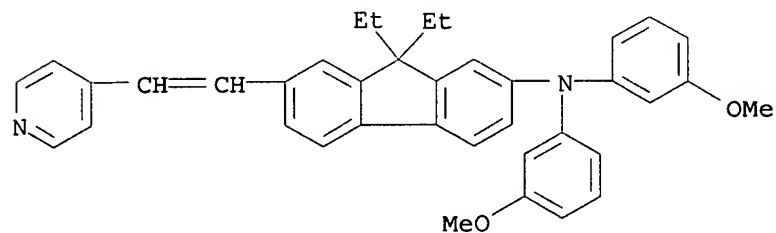
RN 209603-48-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(3-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



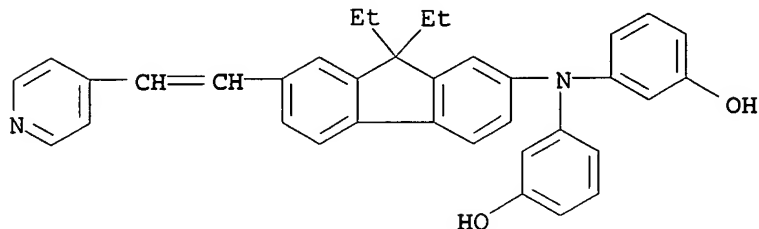
RN 209603-50-9 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-dihexyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



RN 209603-53-2 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-bis(3-methoxyphenyl)-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



RN 209603-56-5 CAPLUS  
 CN Phenol, 3,3'-[[9,9-diethyl-7-[2-(4-pyridinyl)ethenyl]-9H-fluoren-2-yl]imino]bis- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 63 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:33766 CAPLUS

DOCUMENT NUMBER: 128:153818

TITLE: The design and synthesis of new organic molecules with large two-photon absorption cross-sections for optical limiting applications

AUTHOR(S): Reinhardt, B. A.; Brott L. L.; Clarson, S. J.; Kannan, R.; Dillard, A. G.

CORPORATE SOURCE: U.S. Air Force Wright Laboratory, Polymer Branch, WL/MLBP Materials Directorate, Wright-Patterson AFB, OH, 45433-7750, USA

SOURCE: Materials Research Society Symposium Proceedings (1997), 479 (Materials for Optical Limiting II), 3-8  
 CODEN: MRSPDH; ISSN: 0272-9172

PUBLISHER: Materials Research Society

DOCUMENT TYPE: Journal

LANGUAGE: English

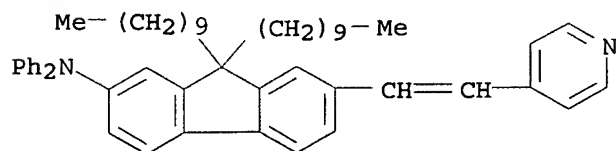
AB The mol. structure/nonlinear optical (NLO) property relationship is explored with seven recently synthesized chromophores. Two sym. compds. were made using electron withdrawing groups separated by an electron rich core while five asym. mols. were developed using electron donating and withdrawing groups coupled by a  $\pi$  electron bridging group. Pendant chains were added to some of the chromophores to improve processability. Their syntheses are described and their optical limiting properties discussed.

IT 191667-13-7P 197969-56-5P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
 (design and preparation of organic mols. with large two-photon absorption cross-sections for optical limiting applications)

RN 191667-13-7 CAPLUS

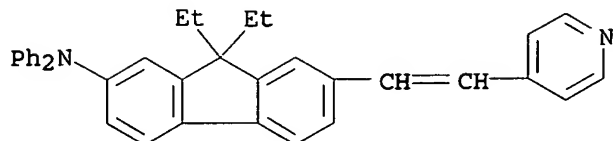
CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]- (9CI) (CA INDEX NAME)



RN 197969-56-5 CAPLUS



CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-(9CI) (CA INDEX NAME)



REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 64 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1997:702997 CAPLUS

DOCUMENT NUMBER: 127:332778

TITLE: Optical power limiting in solution via two-photon absorption: new aromatic heterocyclic dyes with greatly improved performance

AUTHOR(S): Reinhardt, Bruce A.; Brott, Lawrence L.; Clarson, Stephen J.; Kannan, Ramamurthi; Dillard, Ann G.

CORPORATE SOURCE: Polymer Branch, WL/MLBP Materials Directorate, U. S. Air Force Research Laboratory, Wright-Patterson AFB, OH, 45433-7750, USA

SOURCE: Proceedings of SPIE-The International Society for Optical Engineering (1997), 3146(Nonlinear Optical Liquids and Power Limiters), 2-11  
CODEN: PSISDG; ISSN: 0277-786X

PUBLISHER: SPIE-The International Society for Optical Engineering

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Organic compds. which exhibit optical power limiting exclusively via a two-photon absorption mechanism have shown only little promise for providing the limiting activity necessary for the practical protection of eyes and sensors. Unfortunately, there have been few systematic studies of the mol. structure/two-photon absorption property relationships for orgs. documented in the literature. In order to enable the design and synthesis of new mols. with much larger two-photon absorption cross-sections and improved limiting properties, the synthetic chemist must have access to well defined structure/property data. In an attempt to fill this void, work has centered on the design and synthesis of several new families of aromatic heterocyclic chromophores with systematic variations in their mol. structures. Careful characterization of these new materials in solution has produced some well-defined structure/two-photon property relationships at 800 nm. The design and synthesis of these materials are discussed with special emphasis of how the flexibility of the synthetic scheme employed enables the incorporation of these chromophores into a wide variety of materials forms. The characterization of the two-photon properties of these materials and the relationship of these results to their optical limiting behavior in solution will also be reviewed.

IT 191667-13-7P 197969-56-5P

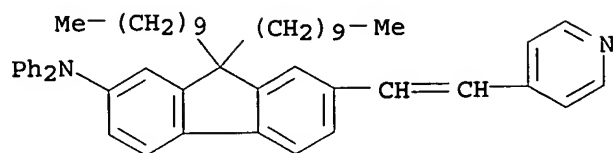
RL: PRP (Properties); SPN (Synthetic preparation); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)

(preparation of aromatic heterocyclic dyes with optical power limiting in solution via two-photon absorption)

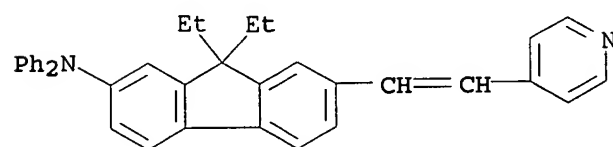
RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-

(9CI) (CA INDEX NAME)



RN 197969-56-5 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-diethyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
 (9CI) (CA INDEX NAME)



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 65 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1997:311012 CAPLUS

DOCUMENT NUMBER: 127:87642

TITLE: Nonlinear optical properties of a new chromophore  
 AUTHOR(S): He, Guang S.; Yuan, Lixiang; Cheng, Ning; Bhawalkar, Jayant D.; Prasad, Paras N.; Brott, Lawrence L.; Clarson, Stephen J.; Reinhardt, Bruce A.

CORPORATE SOURCE: Photonics Research Laboratory, Department Chemistry, State University New York, Buffalo, NY, 14260-3000, USA

SOURCE: Journal of the Optical Society of America B: Optical Physics (1997), 14(5), 1079-1087  
 CODEN: JOBPDE; ISSN: 0740-3224

PUBLISHER: Optical Society of America

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Strong two-photon-absorption (TPA) based nonlinear optical properties of a new chromophore, N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-9,9-di-n-decyl-fluoren-2-amine (AF-50) was exptl. studied. Under excitation with 10-Hz, .apprx.8-ns, and .apprx.800-nm laser pulses, the TPA cross section and the TPA-induced frequency-upconverted emission spectra are measured for AF-50 solns. in various solvents. The most attractive feature of this chromophore is its remarkably high value of the mol. TPA cross section (.apprx.78 + 10-20 cm2/GW in benzene solution). Based on this feature, superior optical power limiting and stabilization performance was demonstrated in a 1-cm-long AF-50 solution sample with concentration of d0 = 0.045

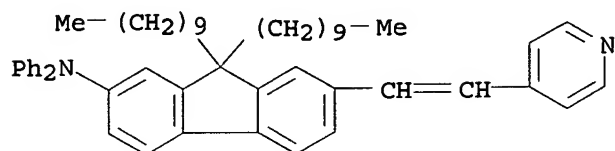
M/L. The nonlinear transmission of the measured sample decreased from .apprx.0.93 to .apprx.0.3 when the input-beam intensity increased from .apprx.10 MW/cm2; the relative intensity fluctuation of the output laser pulses was reduced to 1/3 of that of the input laser beam.

IT 191667-13-7

RL: PRP (Properties)  
 (nonlinear and linear optical properties of new chromophore)

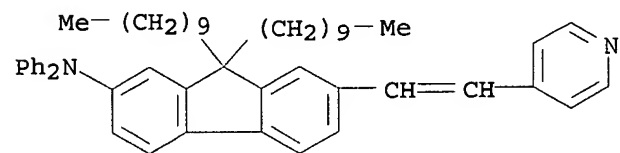
RN 191667-13-7 CAPLUS

CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



REFERENCE COUNT: 26 THERE ARE 26 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

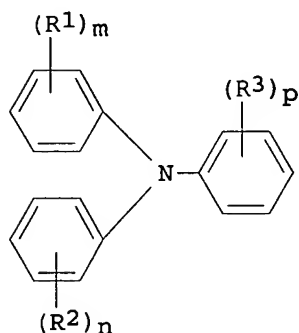
L6 ANSWER 66 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1997:213235 CAPLUS  
 DOCUMENT NUMBER: 126:349333  
 TITLE: Measurement of two-photon absorption cross sections of dye molecules doped in thin films of polymethylmethacrylate  
 AUTHOR(S): Mukherjee, N.; Mukherjee, A.; Reinhardt, B. A.  
 CORPORATE SOURCE: Laser-Matter Interaction Labs Incorporated, Albuquerque, NM, 87110, USA  
 SOURCE: Applied Physics Letters (1997), 70(12), 1524-1526  
 CODEN: APPLAB; ISSN: 0003-6951  
 PUBLISHER: American Institute of Physics  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 AB Two-photon absorption cross sections of 2 dye mols., doped in thin films of polymethylmethacrylate, are measured by comparing fluorescence intensities induced by single- and 2-photon excitations. Ratio of the single- and 2-photon absorption cross sections is related to the ratio of the single- and 2-photon pumped fluorescence intensities. Using this sensitive and background-free detection technique, dispersion of the nonlinearity is measured over a wavelength range of 700-800 nm.  
 IT 191667-13-7, AF 50  
 RL: PRP (Properties)  
 (measurement of two-photon absorption cross sections of dye mols. doped in thin films of polymethylmethacrylate)  
 RN 191667-13-7 CAPLUS  
 CN 9H-Fluoren-2-amine, 9,9-didecyl-N,N-diphenyl-7-[2-(4-pyridinyl)ethenyl]-  
(9CI) (CA INDEX NAME)



L6 ANSWER 67 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1995:528442 CAPLUS  
 DOCUMENT NUMBER: 122:277533  
 TITLE: Organic electroluminescence device containing carbonyl tryarylamine compounds  
 INVENTOR(S): Kikuchi, Norihiro; Nakano, Takashi  
 PATENT ASSIGNEE(S): Canon Kk, Japan  
 SOURCE: Jpn. Kokai Tokkyo Koho, 19 pp.

DOCUMENT TYPE: CODEN: JKXXAF  
 LANGUAGE: Patent  
 FAMILY ACC. NUM. COUNT: 2 Japanese  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 06256759	A2	19940913	JP 1992-134527	19920428
JP 2939052	B2	19990825		
US 5378519	A	19950103	US 1993-52460	19930426
PRIORITY APPLN. INFO.:			JP 1992-134526	A 19920428
			JP 1992-134527	A 19920428
OTHER SOURCE(S):	MARPAT 122:277533			
GI				



AB An electroluminescence device comprising a cathode, an anode, and organic layer(s) formed between the electrodes, is characterized in that the one of the organic layers comprises a CO-containing amine derivative I (R1, R2, R3

=

alkyl, aralkyl, aromatic, heterocyclic, alkoxy, aryloxy, halo, nitro, cyano, hydroxyl, and amino; m, n, p, = 0-5).

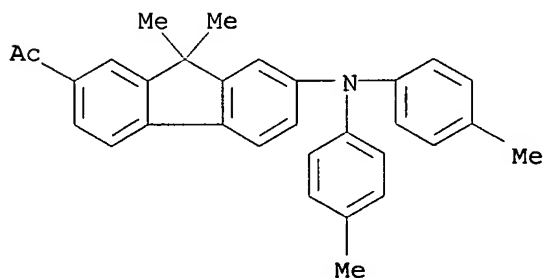
IT 162815-83-0 162815-86-3

RL: DEV (Device component use); USES (Uses)

(organic electroluminescence device containing carbonyl tryarylamine compds.)

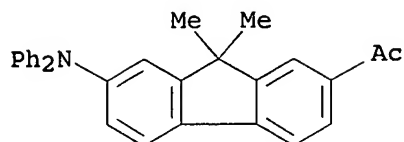
RN 162815-83-0 CAPLUS

CN Ethanone, 1-[7-[bis(4-methylphenyl)amino]-9,9-dimethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



RN 162815-86-3 CAPLUS

CN Ethanone, 1-[7-(diphenylamino)-9,9-dimethyl-9H-fluoren-2-yl]- (9CI) (CA INDEX NAME)



L6 ANSWER 68 OF 68 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1987:565420 CAPLUS

DOCUMENT NUMBER: 107:165420

TITLE: Electrophotographic charge-generating tetrakisazo pigments

INVENTOR(S): Matsumoto, Masakazu; Umehara, Masashige; Takiguchi, Takao; Yamashita, Masataka; Ishikawa, Shozo

PATENT ASSIGNEE(S): Canon K. K., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 40 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent

LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 6

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 62018565	A2	19870127	JP 1985-157699	19850717
JP 04035750	B4	19920612		
US 4666810	A	19870519	US 1986-852243	19860415
PRIORITY APPLN. INFO.:			JP 1985-80248	A 19850417
			JP 1985-157699	A 19850717
			JP 1985-157700	A 19850717
			JP 1985-159401	A 19850718
			JP 1985-159402	A 19850718
			JP 1985-159403	A 19850718

AB The charge-generating tetrakisazo pigments have the formula (AN:NZ3)(AN:NZ4)NZ1CB1:CB2Z2N(Z5N:NA)(Z6N:NA) (I; A = coupler residue with a phenolic OH group; Z1-Z6 = arylene, condensed polycyclene, heterocyclene; B1, B2 = H, halo, CF3, CN, etc.). An electrophotog. charge-generating layer may contain a tetrakisazo pigment of the formula I (A = coupler residue from 3-hydroxy-2-naphthoic acid anilide; Z1-Z6 = 1,4-phenylene; B1, B2 = H) and a poly(vinyl butyral) binder. It provides electrophotog. photoreceptors with improved sensitivity and voltage stability for repeated use.

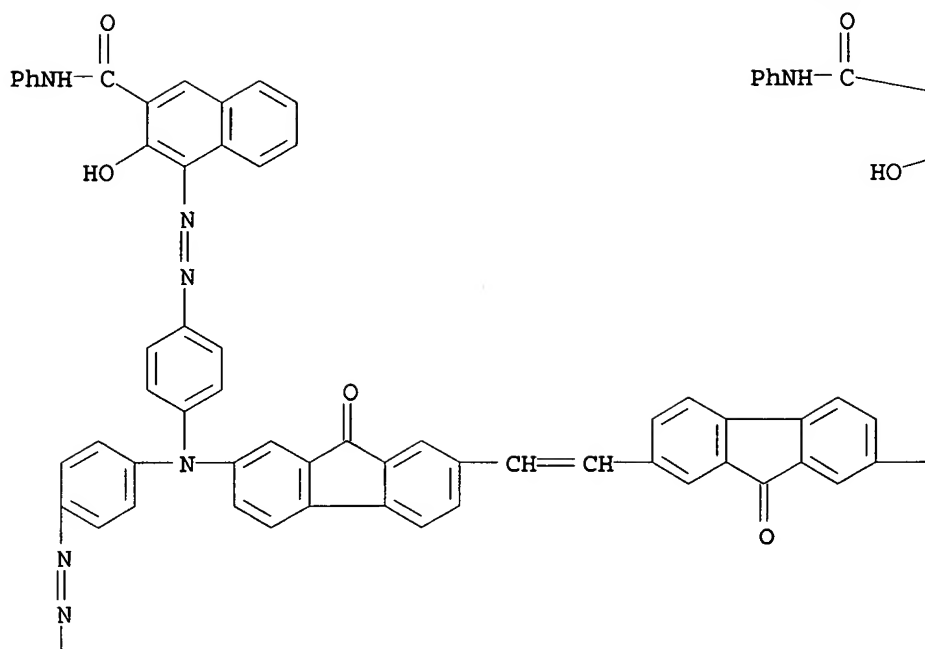
IT 110573-50-7 110573-52-9 110573-72-3

RL: USES (Uses)

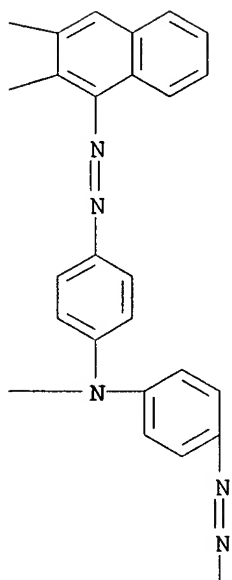
(electrophotog. charge-generating pigments)

RN 110573-50-7 CAPLUS

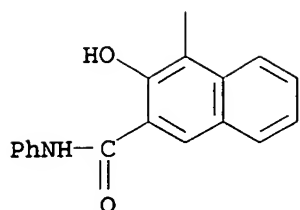
CN 2-Naphthalenecarboxamide, 4,4',4'',4'''-[1,2-ethenediylbis[(9-oxo-9H-fluorene-7,2-diyl)nitrilobis(4,1-phenyleneazo)]]tetrakis[3-hydroxy-N-phenyl]- (9CI) (CA INDEX NAME)



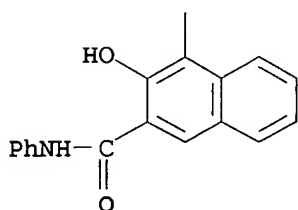
PAGE 1-B



PAGE 2-A



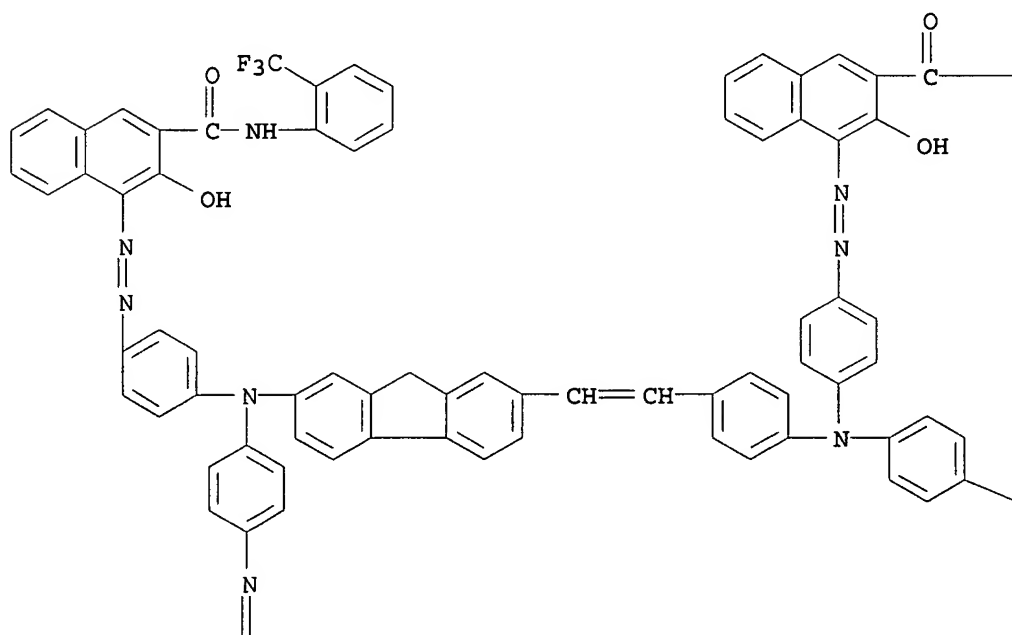
PAGE 2-B

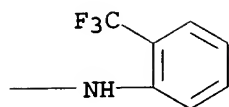


RN 110573-52-9 CAPLUS

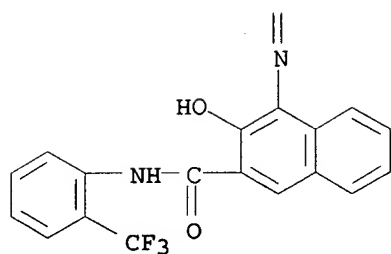
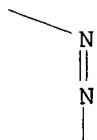
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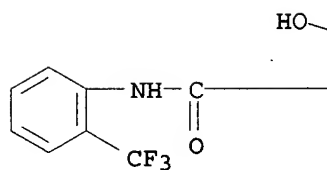




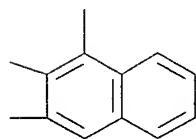
PAGE 1-B



PAGE 2-A



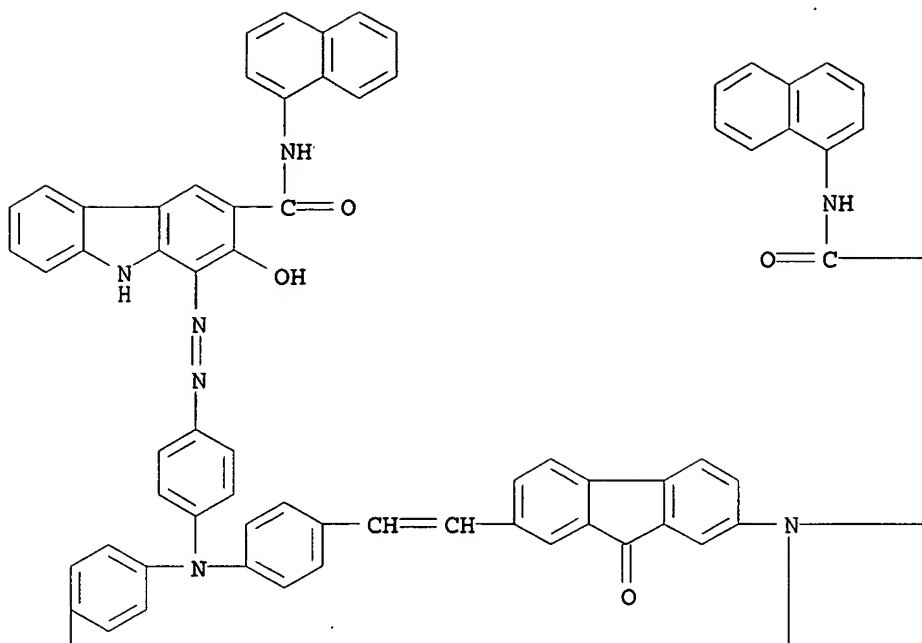
PAGE 2-B



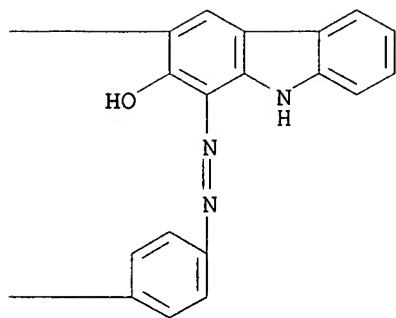
RN 110573-72-3 CAPLUS  
 CN 9H-Carbazole-3-carboxamide, 1,1'-[[[4-[2-[7-[bis[4-[[2-hydroxy-3-[(1-naphthalenylamino) carbonyl]-9H-carbazol-1-yl]azo]phenyl]amino]-9-oxo-9H-fluoren-2-yl]ethenyl]phenyl]imino]bis(4,1-phenyleneazo)]bis[2-hydroxy-N-1-naphthalenyl- (9CI) (CA INDEX NAME)

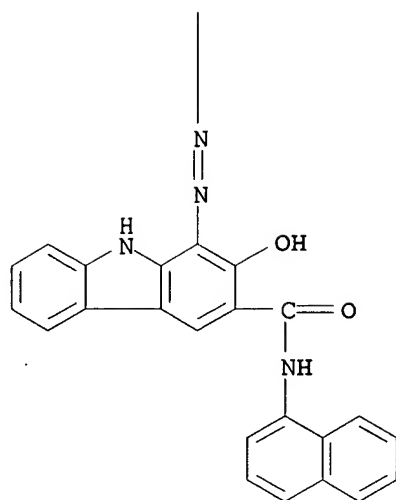


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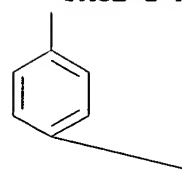


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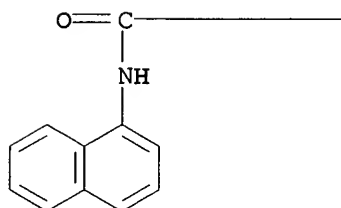




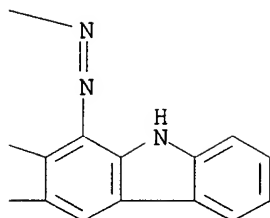
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COST IN U.S. DOLLARS

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

CA SUBSCRIBER PRICE

SINCE FILE

ENTRY

349.81

SINCE FILE

ENTRY

-51.00

TOTAL

SESSION

518.78

TOTAL

SESSION

-51.00

STN INTERNATIONAL LOGOFF AT 13:36:25 ON 19 JUN 2006